



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 08:25 AM GMT

PDB ID : 2R4L
Title : Crystal structure of the long-chain fatty acid transporter FadL mutant P34A
Authors : Hearn, E.M.; Patel, D.R.; Lepore, B.W.; Indic, M.; van den Berg, B.
Deposited on : 2007-08-31
Resolution : 3.30 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>

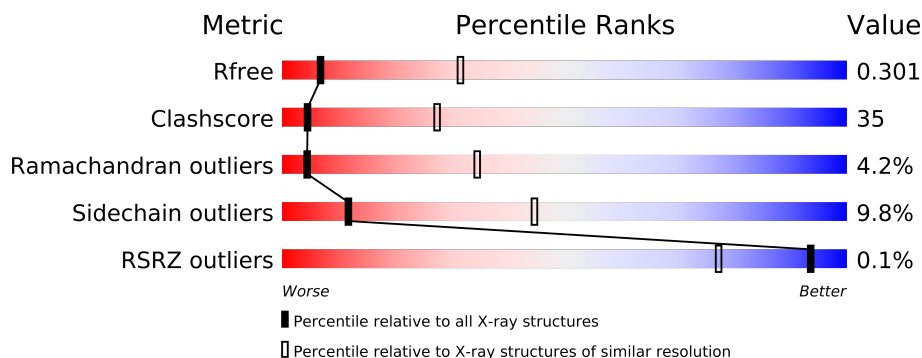
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.30 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	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (3.40-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	427	
1	B	427	
1	C	427	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	LDA	A	505	-	X
2	LDA	A	506	-	X
2	LDA	B	503	-	X
2	LDA	B	504	-	X
2	LDA	C	501	-	X
2	LDA	C	502	-	X

2 Entry composition ⓘ

There are 2 unique types of molecules in this entry. The entry contains 9740 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 Long-chain fatty acid transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	412	Total	C	N	O	S	0	0	0
			3184	2016	539	623	6			
1	B	421	Total	C	N	O	S	0	0	0
			3249	2053	552	638	6			
1	C	416	Total	C	N	O	S	0	0	0
			3211	2032	544	629	6			

There are 24 discrepancies between the modelled and reference sequences:

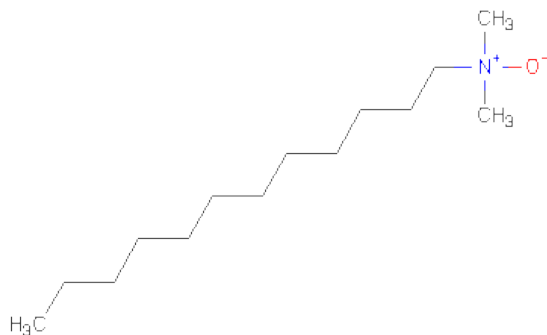
Chain	Residue	Modelled	Actual	Comment	Reference
A	34	ALA	PRO	ENGINEERED	UNP P10384
A	197	THR	ILE	CONFLICT	UNP P10384
A	422	HIS	-	EXPRESSION TAG	UNP P10384
A	423	HIS	-	EXPRESSION TAG	UNP P10384
A	424	HIS	-	EXPRESSION TAG	UNP P10384
A	425	HIS	-	EXPRESSION TAG	UNP P10384
A	426	HIS	-	EXPRESSION TAG	UNP P10384
A	427	HIS	-	EXPRESSION TAG	UNP P10384
B	34	ALA	PRO	ENGINEERED	UNP P10384
B	197	THR	ILE	CONFLICT	UNP P10384
B	422	HIS	-	EXPRESSION TAG	UNP P10384
B	423	HIS	-	EXPRESSION TAG	UNP P10384
B	424	HIS	-	EXPRESSION TAG	UNP P10384
B	425	HIS	-	EXPRESSION TAG	UNP P10384
B	426	HIS	-	EXPRESSION TAG	UNP P10384
B	427	HIS	-	EXPRESSION TAG	UNP P10384
C	34	ALA	PRO	ENGINEERED	UNP P10384
C	197	THR	ILE	CONFLICT	UNP P10384
C	422	HIS	-	EXPRESSION TAG	UNP P10384
C	423	HIS	-	EXPRESSION TAG	UNP P10384
C	424	HIS	-	EXPRESSION TAG	UNP P10384
C	425	HIS	-	EXPRESSION TAG	UNP P10384
C	426	HIS	-	EXPRESSION TAG	UNP P10384

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Chain	Residue	Modelled	Actual	Comment	Reference
C	427	HIS	-	EXPRESSION TAG	UNP P10384

- Molecule 2 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: $C_{14}H_{31}NO$).



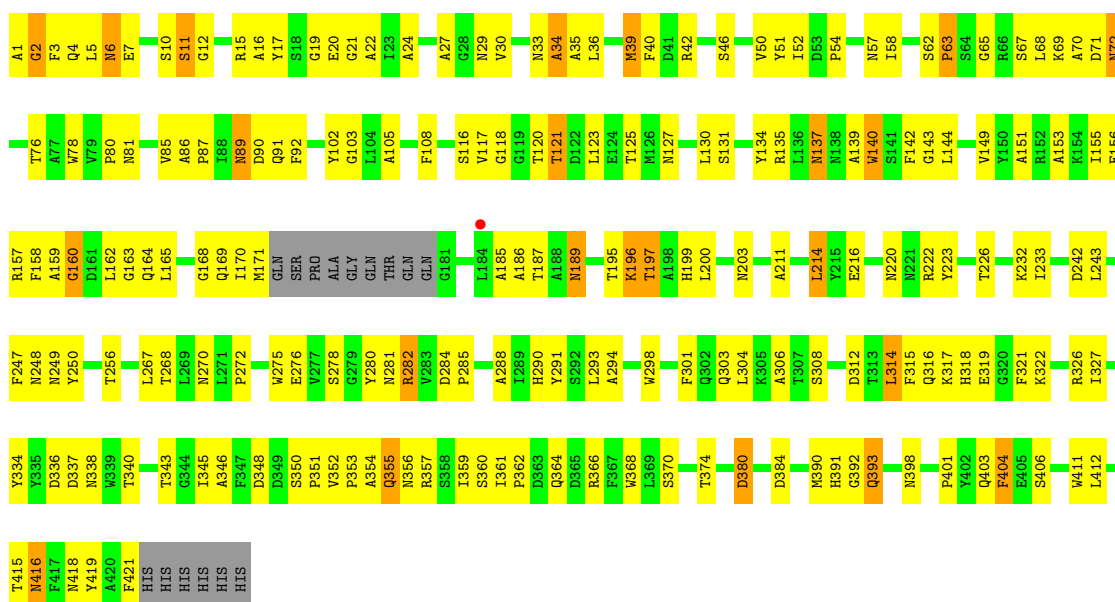
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			16	14	1	1		
2	C	1	Total	C	N	O	0	0
			16	14	1	1		
2	B	1	Total	C	N	O	0	0
			16	14	1	1		
2	B	1	Total	C	N	O	0	0
			16	14	1	1		
2	A	1	Total	C	N	O	0	0
			16	14	1	1		
2	A	1	Total	C	N	O	0	0
			16	14	1	1		

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.

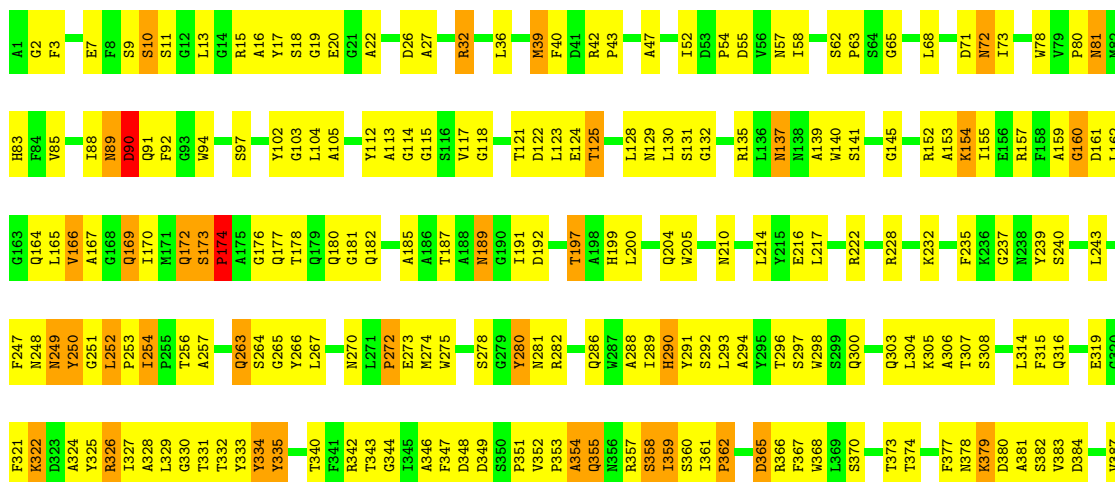
• Molecule 1: Long-chain fatty acid transport protein

Chain A:



• Molecule 1: Long-chain fatty acid transport protein

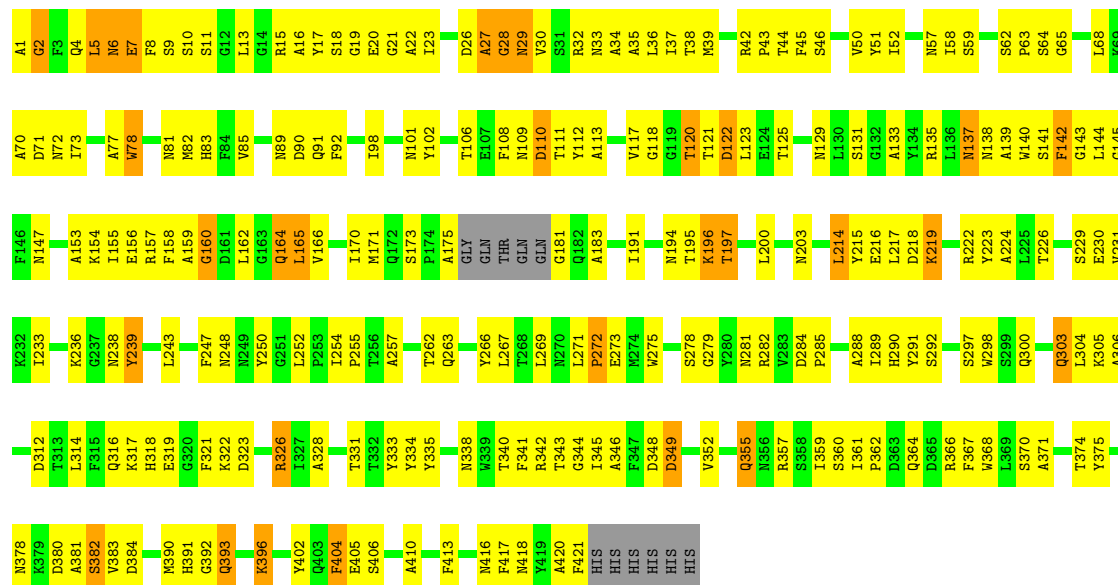
Chain B:





• Molecule 1: Long-chain fatty acid transport protein

Chain C:



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	112.80Å 167.04Å 197.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 3.30 48.40 – 3.30	Depositor EDS
% Data completeness (in resolution range)	95.2 (10.00-3.30) 94.8 (48.40-3.30)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.92 (at 3.33Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.302 0.240 , 0.301	Depositor DCC
R_{free} test set	2602 reflections (10.01%)	DCC
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.683	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , -10.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 27032 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	9740	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.49	0/3270	0.74	0/4450
1	B	0.47	0/3337	0.71	0/4543
1	C	0.47	0/3298	0.74	0/4489
All	All	0.48	0/9905	0.73	0/13482

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	291	TYR	Sidechain
1	C	239	TYR	Sidechain

5.2 Close contacts ⓘ

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

and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3184	0	2967	194	0
1	B	3249	0	3027	222	0
1	C	3211	0	2992	254	0
2	A	32	0	62	18	0
2	B	32	0	62	11	0
2	C	32	0	62	19	0
All	All	9740	0	9172	668	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 35.

All (668) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:34:ALA:O	1:C:143:GLY:HA3	1.46	1.15
1:C:361:ILE:HD11	2:C:501:LDA:H22	1.22	1.10
1:C:58:ILE:HB	1:C:70:ALA:HB3	1.39	1.05
1:B:346:ALA:HB3	1:B:368:TRP:HB2	1.40	1.03
1:B:252:LEU:HB3	1:B:254:ILE:HG12	1.41	1.00
1:B:382:SER:HB3	1:B:418:ASN:HB2	1.45	0.96
1:C:64:SER:HB3	1:C:165:LEU:HD21	1.49	0.94
1:C:10:SER:H	1:C:416:ASN:HD22	1.18	0.92
1:A:293:LEU:HD13	1:A:327:ILE:HG12	1.49	0.92
1:A:10:SER:OG	1:A:416:ASN:HB2	1.70	0.91
1:C:21:GLY:HA2	1:C:36:LEU:HD11	1.50	0.90
1:C:10:SER:H	1:C:416:ASN:ND2	1.68	0.90
1:A:21:GLY:H	1:A:29:ASN:HD22	1.17	0.89
1:A:117:VAL:HG23	1:A:359:ILE:HD11	1.55	0.88
1:B:293:LEU:HD13	1:B:327:ILE:CD1	2.04	0.87
1:C:368:TRP:HZ2	2:C:502:LDA:HM23	1.39	0.87
1:C:303:GLN:HE22	1:C:316:GLN:HE21	1.22	0.87
1:B:63:PRO:HA	1:C:138:ASN:OD1	1.74	0.87
1:B:293:LEU:HD13	1:B:327:ILE:HD11	1.58	0.86
1:C:278:SER:HB3	1:C:292:SER:HB3	1.56	0.85
1:A:319:GLU:OE2	2:A:506:LDA:HM22	1.76	0.85
1:C:346:ALA:HB3	1:C:368:TRP:HB2	1.58	0.84
1:A:102:TYR:O	1:A:125:THR:HB	1.78	0.83
1:C:390:MET:HE1	2:C:502:LDA:HM22	1.61	0.83
1:A:354:ALA:O	1:A:355:GLN:HB2	1.78	0.82
1:A:343:THR:HG23	1:A:370:SER:O	1.80	0.82

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:267:LEU:HD23	1:A:268:THR:N	1.95	0.81
1:C:390:MET:HG3	1:C:410:ALA:HB3	1.60	0.81
1:C:361:ILE:HD11	2:C:501:LDA:C2	2.10	0.80
1:B:319:GLU:OE1	2:B:503:LDA:HM21	1.82	0.80
1:C:110:ASP:O	1:C:194:ASN:HB3	1.82	0.79
1:C:38:THR:HG21	1:C:141:SER:OG	1.81	0.79
1:C:21:GLY:H	1:C:29:ASN:HD22	1.27	0.79
1:C:131:SER:HB3	1:C:145:GLY:HA3	1.63	0.79
1:C:27:ALA:HB2	1:C:44:THR:HG22	1.63	0.79
1:A:89:ASN:ND2	1:A:90:ASP:H	1.81	0.79
1:B:306:ALA:HB3	1:B:315:PHE:HB3	1.63	0.78
1:A:196:LYS:HB2	1:A:242:ASP:OD2	1.84	0.78
1:A:164:GLN:HG3	1:A:189:ASN:ND2	1.99	0.78
1:B:102:TYR:O	1:B:125:THR:HB	1.85	0.77
1:C:281:ASN:HB2	1:C:289:ILE:HG23	1.66	0.76
1:A:4:GLN:C	1:A:5:LEU:HD12	2.06	0.76
1:A:153:ALA:CB	2:A:506:LDA:H91	2.16	0.76
1:C:157:ARG:HH12	2:C:501:LDA:HM22	1.51	0.75
1:B:62:SER:HB2	1:B:68:LEU:HD21	1.67	0.75
1:B:131:SER:HB3	1:B:145:GLY:HA2	1.69	0.75
1:A:200:LEU:HB3	2:A:506:LDA:H123	1.67	0.74
1:A:303:GLN:HE22	1:A:316:GLN:HE21	1.33	0.74
1:B:199:HIS:C	1:B:200:LEU:HD12	2.08	0.74
1:A:168:GLY:O	1:A:171:MET:HG2	1.88	0.74
1:C:137:ASN:HD21	1:C:140:TRP:H	1.34	0.73
1:C:164:GLN:O	1:C:166:VAL:N	2.20	0.73
1:C:137:ASN:HD21	1:C:140:TRP:N	1.86	0.73
1:A:276:GLU:HG3	1:A:294:ALA:HB2	1.70	0.73
1:C:131:SER:HB3	1:C:145:GLY:CA	2.18	0.73
1:C:278:SER:HB3	1:C:292:SER:CB	2.19	0.72
1:B:154:LYS:HA	1:B:200:LEU:O	1.88	0.72
1:C:137:ASN:HD22	1:C:137:ASN:C	1.92	0.72
1:C:281:ASN:ND2	1:C:291:TYR:HE2	1.88	0.72
1:B:19:GLY:H	1:B:292:SER:HB3	1.55	0.72
1:C:338:ASN:HB3	1:C:375:TYR:HE1	1.55	0.72
1:B:273:GLU:H	1:B:300:GLN:NE2	1.88	0.72
1:C:102:TYR:O	1:C:125:THR:HB	1.89	0.71
1:B:42:ARG:HD2	1:B:421:PHE:O	1.91	0.71
1:C:254:ILE:HG13	1:C:255:PRO:HD2	1.72	0.71
1:B:65:GLY:C	1:C:91:GLN:HE21	1.94	0.71
1:B:396:LYS:HD2	1:B:396:LYS:H	1.56	0.70
1:B:396:LYS:N	1:B:396:LYS:HD2	2.06	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:393:GLN:HE21	1:B:393:GLN:HA	1.55	0.70
1:A:17:TYR:HA	1:A:20:GLU:OE1	1.90	0.70
1:A:5:LEU:HA	2:A:505:LDA:HM13	1.72	0.70
1:C:5:LEU:HD22	1:C:6:ASN:H	1.57	0.70
1:A:364:GLN:HG3	1:A:393:GLN:O	1.91	0.69
1:B:346:ALA:HB3	1:B:368:TRP:CB	2.20	0.69
1:C:20:GLU:HA	1:C:20:GLU:OE2	1.92	0.69
1:A:232:LYS:HE2	1:A:270:ASN:ND2	2.07	0.69
1:B:303:GLN:HE22	1:B:316:GLN:NE2	1.90	0.69
1:C:21:GLY:CA	1:C:36:LEU:HD11	2.20	0.69
1:A:352:VAL:HG11	1:A:357:ARG:HA	1.75	0.69
1:A:415:THR:HG22	1:A:416:ASN:H	1.57	0.68
1:B:178:THR:HG22	1:B:180:GLN:H	1.56	0.68
1:C:374:THR:HG23	1:C:384:ASP:OD2	1.94	0.68
1:C:52:ILE:HG21	2:C:502:LDA:H52	1.76	0.68
1:B:378:ASN:O	1:B:379:LYS:HB2	1.94	0.68
1:C:393:GLN:HA	1:C:393:GLN:HE21	1.58	0.68
1:C:27:ALA:O	1:C:30:VAL:HG22	1.94	0.67
1:C:7:GLU:HG2	1:C:50:VAL:HG11	1.76	0.67
1:B:131:SER:HB3	1:B:145:GLY:CA	2.24	0.67
1:C:303:GLN:HE22	1:C:316:GLN:NE2	1.92	0.66
1:B:135:ARG:HG3	1:B:137:ASN:O	1.95	0.66
1:C:338:ASN:HB3	1:C:375:TYR:CE1	2.31	0.66
1:C:281:ASN:HB2	1:C:289:ILE:CG2	2.26	0.66
1:B:199:HIS:O	1:B:200:LEU:HD12	1.96	0.66
1:A:33:ASN:HB2	1:A:226:THR:HG21	1.78	0.66
1:B:272:PRO:HA	1:B:300:GLN:HE21	1.59	0.65
1:B:17:TYR:HA	1:B:20:GLU:OE1	1.96	0.65
1:A:89:ASN:HD22	1:A:90:ASP:H	1.44	0.65
1:B:306:ALA:O	1:B:314:LEU:HB2	1.97	0.65
1:C:7:GLU:HA	1:C:17:TYR:OH	1.96	0.65
1:B:10:SER:HB3	1:B:416:ASN:ND2	2.12	0.65
1:A:326:ARG:HG3	1:A:348:ASP:HB3	1.76	0.65
1:C:305:LYS:HD2	1:C:316:GLN:NE2	2.12	0.65
1:B:68:LEU:HD13	1:B:404:PHE:CE2	2.32	0.65
1:B:137:ASN:HD21	1:B:140:TRP:H	1.44	0.65
1:B:170:ILE:HG21	1:B:185:ALA:HB2	1.79	0.65
1:C:346:ALA:HB3	1:C:368:TRP:CB	2.28	0.64
1:A:67:SER:C	1:A:68:LEU:HD23	2.18	0.64
1:C:173:SER:O	1:C:175:ALA:N	2.31	0.64
1:C:110:ASP:O	1:C:194:ASN:CB	2.45	0.64
1:B:237:GLY:O	1:B:265:GLY:N	2.30	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:157:ARG:HB2	1:C:197:THR:HG22	1.80	0.63
1:C:380:ASP:HA	1:C:420:ALA:HB3	1.81	0.63
1:A:284:ASP:OD1	1:A:285:PRO:HD2	1.98	0.63
1:A:7:GLU:OE1	1:A:7:GLU:N	2.32	0.63
1:C:6:ASN:O	1:C:7:GLU:HB2	1.98	0.63
1:B:293:LEU:HD13	1:B:327:ILE:HD13	1.80	0.62
1:B:52:ILE:O	1:B:54:PRO:HD3	1.99	0.62
1:A:360:SER:C	1:A:362:PRO:HD3	2.20	0.62
1:C:238:ASN:HD22	1:C:262:THR:HG21	1.63	0.62
1:C:396:LYS:H	1:C:396:LYS:HD2	1.65	0.62
1:C:137:ASN:C	1:C:137:ASN:ND2	2.52	0.62
1:C:321:PHE:CE2	1:C:352:VAL:HG22	2.34	0.62
1:A:303:GLN:HE22	1:A:316:GLN:NE2	1.95	0.62
1:B:137:ASN:ND2	1:B:139:ALA:H	1.96	0.62
1:A:415:THR:HG22	1:A:416:ASN:N	2.13	0.62
1:B:32:ARG:HH11	1:B:32:ARG:HG2	1.65	0.62
1:A:42:ARG:HD2	1:A:421:PHE:O	1.99	0.62
1:C:10:SER:N	1:C:416:ASN:HD22	1.94	0.61
1:C:106:THR:HG21	1:C:360:SER:HA	1.81	0.61
1:C:272:PRO:HA	1:C:300:GLN:HE21	1.66	0.61
1:C:303:GLN:NE2	1:C:316:GLN:HE21	1.94	0.61
1:C:159:ALA:O	1:C:160:GLY:C	2.38	0.61
1:A:92:PHE:CE2	1:A:134:TYR:HB2	2.36	0.60
1:A:398:ASN:OD1	1:A:403:GLN:HG3	2.00	0.60
1:A:360:SER:O	1:A:362:PRO:HD3	2.02	0.60
1:C:352:VAL:HG11	1:C:357:ARG:HA	1.84	0.60
1:B:228:ARG:HB3	1:B:274:MET:HB3	1.84	0.60
1:B:57:ASN:HD22	1:B:71:ASP:HA	1.67	0.60
1:C:26:ASP:HB2	1:C:418:ASN:ND2	2.17	0.59
1:C:19:GLY:H	1:C:292:SER:CB	2.15	0.59
1:A:308:SER:HB3	1:A:314:LEU:HD11	1.83	0.59
1:C:51:TYR:N	1:C:413:PHE:O	2.21	0.59
1:A:304:LEU:HD11	2:A:506:LDA:H32	1.85	0.59
1:A:345:ILE:HA	1:A:368:TRP:O	2.02	0.59
1:B:358:SER:HB2	1:B:399:GLU:CD	2.22	0.59
1:C:18:SER:OG	1:C:290:HIS:HD2	1.85	0.59
1:B:10:SER:CB	1:B:416:ASN:HD22	2.16	0.59
1:A:243:LEU:HD22	1:A:247:PHE:CE1	2.38	0.59
1:A:12:GLY:HA2	1:A:15:ARG:CZ	2.33	0.59
1:C:38:THR:HG21	1:C:141:SER:HG	1.68	0.59
1:B:89:ASN:O	1:B:91:GLN:N	2.35	0.59
1:B:281:ASN:HB2	1:B:289:ILE:HG23	1.84	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:319:GLU:HB3	1:A:321:PHE:CE2	2.37	0.58
1:B:396:LYS:HG3	1:B:405:GLU:OE2	2.04	0.58
1:C:155:ILE:HD11	2:C:501:LDA:H62	1.86	0.58
1:C:378:ASN:OD1	1:C:380:ASP:N	2.35	0.58
1:B:347:PHE:HE1	1:B:365:ASP:HB3	1.69	0.58
1:C:236:LYS:HE3	1:C:266:TYR:OH	2.04	0.58
1:C:120:THR:HG23	1:C:122:ASP:OD1	2.03	0.58
1:A:293:LEU:CD1	1:A:327:ILE:HG12	2.29	0.58
1:B:273:GLU:N	1:B:300:GLN:NE2	2.51	0.58
1:A:355:GLN:C	1:A:356:ASN:HD22	2.06	0.58
1:B:55:ASP:HB3	1:B:409:LYS:HG2	1.85	0.58
1:C:29:ASN:HB3	1:C:36:LEU:CD1	2.34	0.58
1:A:62:SER:O	1:A:65:GLY:N	2.33	0.58
1:A:39:MET:HE1	1:A:216:GLU:OE1	2.03	0.58
1:B:170:ILE:HD12	1:B:185:ALA:HB2	1.84	0.58
1:B:157:ARG:HB2	1:B:197:THR:HG22	1.85	0.58
1:C:191:ILE:HD11	1:C:247:PHE:CZ	2.38	0.58
1:B:27:ALA:HA	1:B:40:PHE:CZ	2.39	0.58
1:A:58:ILE:HB	1:A:70:ALA:HB3	1.86	0.58
1:C:239:TYR:CD2	1:C:314:LEU:HD23	2.39	0.58
1:C:223:TYR:HA	1:C:279:GLY:HA2	1.85	0.57
1:A:135:ARG:NH2	1:A:216:GLU:OE2	2.37	0.57
1:C:368:TRP:HZ2	2:C:502:LDA:CM2	2.16	0.57
1:A:195:THR:O	1:A:197:THR:N	2.36	0.57
1:C:35:ALA:C	1:C:37:ILE:H	2.07	0.57
1:B:402:TYR:HB2	1:B:404:PHE:CE1	2.39	0.57
1:B:81:ASN:HD21	1:B:416:ASN:HD21	1.52	0.57
1:C:267:LEU:HD12	1:C:306:ALA:HB2	1.86	0.57
1:A:304:LEU:HD12	2:A:506:LDA:HM21	1.86	0.57
1:A:33:ASN:C	1:A:35:ALA:H	2.08	0.57
1:A:155:ILE:HD11	2:A:506:LDA:H81	1.86	0.57
1:B:26:ASP:O	1:B:40:PHE:HZ	1.88	0.57
1:A:354:ALA:HA	1:A:357:ARG:CG	2.35	0.57
1:A:19:GLY:O	1:A:22:ALA:HB3	2.05	0.57
1:A:7:GLU:HG2	1:A:50:VAL:HG11	1.87	0.57
1:C:281:ASN:ND2	1:C:291:TYR:CE2	2.70	0.57
1:B:324:ALA:HB2	1:B:351:PRO:HG3	1.87	0.57
1:A:27:ALA:HB1	1:A:85:VAL:HG23	1.87	0.57
1:B:27:ALA:HB1	1:B:85:VAL:HG23	1.86	0.56
1:A:153:ALA:HB3	2:A:506:LDA:H91	1.86	0.56
1:A:153:ALA:HB1	2:A:506:LDA:H91	1.85	0.56
1:C:137:ASN:ND2	1:C:139:ALA:N	2.53	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:273:GLU:H	1:C:300:GLN:NE2	2.03	0.56
1:B:57:ASN:ND2	1:B:71:ASP:HA	2.20	0.56
1:A:391:HIS:ND1	1:A:392:GLY:N	2.53	0.56
1:A:411:TRP:C	1:A:412:LEU:HD12	2.26	0.56
1:B:88:ILE:HB	1:B:92:PHE:O	2.05	0.56
1:B:104:LEU:HG	1:B:105:ALA:N	2.20	0.56
1:B:346:ALA:CB	1:B:368:TRP:HB2	2.27	0.56
1:A:21:GLY:H	1:A:29:ASN:ND2	1.96	0.56
1:B:272:PRO:HG2	1:B:298:TRP:CE3	2.41	0.56
1:C:36:LEU:HD23	1:C:214:LEU:HD13	1.88	0.56
1:C:266:TYR:O	1:C:306:ALA:HA	2.05	0.56
1:B:304:LEU:HD13	2:B:503:LDA:H12	1.88	0.56
1:C:254:ILE:HG23	1:C:255:PRO:O	2.06	0.56
1:B:390:MET:SD	2:B:504:LDA:CM2	2.94	0.56
1:C:250:TYR:C	1:C:252:LEU:H	2.09	0.55
1:C:191:ILE:HD11	1:C:247:PHE:HZ	1.72	0.55
1:B:367:PHE:O	1:B:390:MET:HA	2.05	0.55
1:C:166:VAL:O	1:C:170:ILE:HG13	2.07	0.55
1:C:281:ASN:HD22	1:C:291:TYR:HE2	1.54	0.55
1:A:11:SER:HB2	1:A:384:ASP:OD1	2.06	0.55
1:B:170:ILE:CD1	1:B:185:ALA:HB2	2.37	0.55
1:C:360:SER:OG	2:C:501:LDA:HM21	2.07	0.55
1:C:21:GLY:O	1:C:36:LEU:HD21	2.06	0.55
1:C:367:PHE:O	1:C:390:MET:HA	2.06	0.55
1:C:368:TRP:CZ2	2:C:502:LDA:HM23	2.31	0.55
1:A:12:GLY:HA2	1:A:15:ARG:NH1	2.21	0.55
1:B:159:ALA:O	1:B:160:GLY:C	2.44	0.55
1:C:304:LEU:HD11	2:C:501:LDA:H32	1.89	0.55
1:A:89:ASN:ND2	1:A:90:ASP:N	2.52	0.55
1:C:223:TYR:CD2	1:C:279:GLY:HA3	2.41	0.55
1:B:280:TYR:C	1:B:280:TYR:CD2	2.79	0.55
1:C:10:SER:CB	1:C:416:ASN:HD22	2.20	0.55
1:A:5:LEU:HD13	1:A:103:GLY:N	2.21	0.55
1:A:57:ASN:ND2	1:A:71:ASP:HA	2.22	0.55
1:C:142:PHE:N	1:C:142:PHE:CD1	2.74	0.55
1:A:336:ASP:O	1:A:338:ASN:N	2.40	0.55
1:B:13:LEU:HA	1:B:17:TYR:CE1	2.41	0.55
1:B:374:THR:HG23	1:B:384:ASP:OD2	2.07	0.55
1:B:216:GLU:HA	1:B:222:ARG:HA	1.88	0.55
1:C:343:THR:HG23	1:C:370:SER:O	2.07	0.54
1:A:312:ASP:O	1:A:314:LEU:HD13	2.08	0.54
1:B:83:HIS:HA	1:B:97:SER:HB3	1.90	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:358:SER:HB2	1:B:399:GLU:OE1	2.07	0.54
1:B:304:LEU:CD1	2:B:503:LDA:H11	2.38	0.54
1:B:384:ASP:O	1:B:415:THR:HA	2.08	0.54
1:A:90:ASP:O	1:A:91:GLN:HG2	2.07	0.54
1:B:374:THR:HG23	1:B:383:VAL:O	2.08	0.54
1:B:153:ALA:HB3	2:B:503:LDA:H102	1.89	0.54
1:B:58:ILE:CD1	1:B:406:SER:HB2	2.38	0.54
1:A:364:GLN:OE1	1:A:364:GLN:N	2.39	0.54
1:A:361:ILE:HG13	2:A:506:LDA:H22	1.90	0.54
1:A:67:SER:O	1:A:68:LEU:HD23	2.08	0.54
1:B:329:LEU:HD12	1:B:330:GLY:H	1.72	0.54
1:B:303:GLN:HE22	1:B:316:GLN:HE21	1.51	0.54
1:B:16:ALA:O	1:B:326:ARG:NH2	2.41	0.54
1:C:326:ARG:HG3	1:C:326:ARG:HH11	1.73	0.54
1:B:65:GLY:C	1:C:91:GLN:NE2	2.62	0.53
1:A:419:TYR:HE2	1:A:421:PHE:CE1	2.26	0.53
1:C:273:GLU:HG2	1:C:300:GLN:OE1	2.09	0.53
1:C:10:SER:N	1:C:416:ASN:ND2	2.49	0.53
1:C:254:ILE:HG13	1:C:255:PRO:CD	2.36	0.53
1:A:170:ILE:O	1:A:170:ILE:HG22	2.09	0.53
1:A:155:ILE:CD1	2:A:506:LDA:H81	2.38	0.53
1:B:352:VAL:HG11	1:B:357:ARG:HA	1.91	0.53
1:C:319:GLU:OE1	2:C:501:LDA:HM13	2.08	0.53
1:C:342:ARG:NH1	1:C:374:THR:OG1	2.38	0.53
1:A:199:HIS:C	1:A:200:LEU:HD12	2.29	0.53
1:C:238:ASN:HB3	1:C:262:THR:HG23	1.90	0.53
1:A:151:ALA:CB	1:A:233:ILE:HD13	2.39	0.53
1:A:140:TRP:HA	1:A:214:LEU:O	2.09	0.53
1:C:390:MET:O	1:C:410:ALA:N	2.43	0.52
1:A:33:ASN:O	1:A:35:ALA:N	2.38	0.52
1:C:238:ASN:HB3	1:C:262:THR:CG2	2.39	0.52
1:C:272:PRO:HA	1:C:300:GLN:NE2	2.24	0.52
1:B:294:ALA:HB3	1:B:326:ARG:HB3	1.91	0.52
1:A:137:ASN:HD21	1:A:140:TRP:HD1	1.56	0.52
1:A:108:PHE:HB2	1:A:118:GLY:O	2.08	0.52
1:C:18:SER:HA	1:C:292:SER:OG	2.09	0.52
1:B:360:SER:C	1:B:362:PRO:HD3	2.30	0.52
1:B:243:LEU:HD22	1:B:247:PHE:CD1	2.45	0.52
1:A:359:ILE:O	1:A:362:PRO:HG3	2.08	0.52
1:A:276:GLU:HG3	1:A:294:ALA:CB	2.38	0.52
1:B:303:GLN:NE2	1:B:316:GLN:HE21	2.07	0.52
1:B:176:GLY:O	1:B:182:GLN:HG3	2.08	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:281:ASN:O	1:A:288:ALA:HA	2.09	0.52
1:A:361:ILE:HD11	2:A:506:LDA:H31	1.92	0.52
1:B:308:SER:HB3	1:B:314:LEU:HD11	1.92	0.52
1:C:297:SER:HA	1:C:323:ASP:OD2	2.09	0.52
1:B:94:TRP:HA	1:B:132:GLY:HA2	1.92	0.52
1:A:2:GLY:H	1:A:351:PRO:CG	2.23	0.52
1:B:47:ALA:HA	1:B:81:ASN:O	2.09	0.52
1:C:15:ARG:NH1	1:C:20:GLU:OE1	2.43	0.52
1:A:5:LEU:CD1	1:A:103:GLY:N	2.73	0.52
1:B:19:GLY:HA2	1:B:278:SER:HB2	1.92	0.52
1:B:326:ARG:HB2	1:B:348:ASP:HB2	1.92	0.52
1:B:319:GLU:HB3	1:B:321:PHE:CE2	2.44	0.51
1:A:69:LYS:HE2	1:A:71:ASP:OD2	2.10	0.51
1:A:58:ILE:HD13	1:A:406:SER:HB2	1.92	0.51
1:A:355:GLN:C	1:A:356:ASN:ND2	2.63	0.51
1:A:267:LEU:C	1:A:267:LEU:HD23	2.31	0.51
1:A:19:GLY:HA3	1:A:290:HIS:HB2	1.93	0.51
1:B:173:SER:CB	1:B:174:PRO:HD2	2.38	0.51
1:A:5:LEU:HA	2:A:505:LDA:CM1	2.38	0.51
1:C:9:SER:OG	1:C:11:SER:HB3	2.11	0.51
1:B:173:SER:HB2	1:B:174:PRO:HD2	1.91	0.51
1:C:20:GLU:HB2	1:C:32:ARG:HD3	1.93	0.51
1:C:16:ALA:HB1	1:C:346:ALA:HB2	1.92	0.51
1:C:89:ASN:OD1	1:C:92:PHE:HB2	2.10	0.51
1:B:166:VAL:O	1:B:166:VAL:HG12	2.10	0.51
1:B:204:GLN:OE1	1:B:205:TRP:N	2.34	0.51
1:B:248:ASN:OD1	1:B:257:ALA:HB3	2.11	0.51
1:B:326:ARG:HD2	1:B:348:ASP:HB3	1.92	0.51
1:A:140:TRP:CD1	1:A:140:TRP:N	2.78	0.51
1:B:172:GLN:O	1:B:173:SER:HB3	2.11	0.51
1:C:1:ALA:O	1:C:2:GLY:C	2.47	0.51
1:C:29:ASN:HB3	1:C:36:LEU:HD13	1.93	0.51
1:C:390:MET:CE	2:C:502:LDA:HM22	2.39	0.51
1:C:16:ALA:HB1	1:C:346:ALA:CB	2.40	0.50
1:A:2:GLY:H	1:A:351:PRO:HG2	1.76	0.50
1:B:329:LEU:HD12	1:B:330:GLY:N	2.26	0.50
1:B:252:LEU:HB3	1:B:254:ILE:CG1	2.29	0.50
1:C:396:LYS:H	1:C:396:LYS:CD	2.24	0.50
1:C:68:LEU:HD22	1:C:404:PHE:CE2	2.47	0.50
1:C:81:ASN:HD22	1:C:83:HIS:CE1	2.30	0.50
1:C:272:PRO:HG2	1:C:298:TRP:CD2	2.46	0.50
1:A:39:MET:HE3	1:A:216:GLU:OE2	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:224:ALA:N	1:C:278:SER:O	2.42	0.50
1:A:137:ASN:HD22	1:A:137:ASN:H	1.59	0.50
1:A:351:PRO:O	1:A:353:PRO:HD3	2.12	0.50
1:B:18:SER:HB3	1:B:328:ALA:HB1	1.92	0.50
1:C:42:ARG:HD2	1:C:421:PHE:O	2.12	0.50
1:C:51:TYR:HD1	1:C:78:TRP:HD1	1.60	0.50
1:A:137:ASN:HD21	1:A:139:ALA:HB3	1.75	0.50
1:B:239:TYR:CG	1:B:240:SER:N	2.80	0.50
1:C:273:GLU:N	1:C:300:GLN:HE22	2.09	0.50
1:C:82:MET:HE3	1:C:98:ILE:HD12	1.93	0.50
1:C:248:ASN:HD21	1:C:257:ALA:H	1.58	0.50
1:C:123:LEU:CD1	1:C:153:ALA:HB2	2.42	0.49
1:C:108:PHE:HB2	1:C:118:GLY:O	2.11	0.49
1:C:304:LEU:CD1	2:C:501:LDA:H11	2.43	0.49
1:A:355:GLN:O	1:A:356:ASN:ND2	2.41	0.49
1:A:301:PHE:CE2	1:A:319:GLU:HG3	2.47	0.49
1:B:39:MET:HE1	1:B:135:ARG:HH22	1.77	0.49
1:A:220:ASN:HB3	1:A:282:ARG:CB	2.42	0.49
1:C:164:GLN:CG	1:C:165:LEU:H	2.26	0.49
1:C:62:SER:O	1:C:65:GLY:N	2.29	0.49
1:A:354:ALA:HA	1:A:357:ARG:HG2	1.93	0.49
1:C:137:ASN:ND2	1:C:139:ALA:H	2.09	0.49
1:C:291:TYR:N	1:C:291:TYR:CD2	2.79	0.49
1:B:137:ASN:ND2	1:B:140:TRP:H	2.07	0.49
1:B:239:TYR:OH	1:B:256:THR:O	2.18	0.49
1:C:20:GLU:CB	1:C:32:ARG:HD3	2.42	0.49
1:B:354:ALA:O	1:B:355:GLN:HB3	2.13	0.49
1:C:312:ASP:O	1:C:314:LEU:HD13	2.12	0.49
1:C:1:ALA:CB	1:C:4:GLN:HB3	2.43	0.49
1:C:70:ALA:HB1	1:C:73:ILE:CG2	2.42	0.49
1:C:28:GLY:HA2	1:C:83:HIS:CD2	2.48	0.49
1:C:51:TYR:CD1	1:C:78:TRP:HD1	2.30	0.49
1:B:359:ILE:O	1:B:362:PRO:HG3	2.13	0.49
1:A:103:GLY:HA2	1:A:123:LEU:O	2.13	0.48
1:A:158:PHE:HA	1:A:197:THR:HB	1.95	0.48
1:C:223:TYR:HE2	1:C:291:TYR:HH	1.61	0.48
1:C:156:GLU:OE2	1:C:196:LYS:HE2	2.13	0.48
1:C:333:TYR:N	1:C:341:PHE:O	2.39	0.48
1:C:19:GLY:HA2	1:C:278:SER:HB2	1.95	0.48
1:A:121:THR:OG1	1:A:157:ARG:NH2	2.45	0.48
1:C:326:ARG:HG3	1:C:348:ASP:HB3	1.94	0.48
1:B:177:GLN:HE21	1:B:177:GLN:HA	1.79	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:306:ALA:HB3	1:A:315:PHE:HB3	1.94	0.48
1:B:332:THR:HG22	1:B:334:TYR:CE1	2.48	0.48
1:C:157:ARG:CB	1:C:197:THR:HG22	2.43	0.48
1:A:321:PHE:CE2	1:A:352:VAL:HG22	2.49	0.48
1:C:6:ASN:O	1:C:7:GLU:CB	2.60	0.48
1:A:352:VAL:CG1	1:A:357:ARG:HA	2.42	0.48
1:B:170:ILE:CG2	1:B:185:ALA:HB2	2.43	0.48
1:B:307:THR:O	1:B:314:LEU:HD22	2.13	0.48
1:C:33:ASN:HB2	1:C:226:THR:HG21	1.94	0.48
1:A:366:ARG:HD3	1:A:390:MET:HE3	1.94	0.48
1:B:9:SER:O	1:B:11:SER:N	2.40	0.48
1:C:35:ALA:C	1:C:37:ILE:N	2.67	0.48
1:B:137:ASN:HD22	1:B:137:ASN:C	2.17	0.48
1:C:20:GLU:HB2	1:C:32:ARG:CD	2.44	0.48
1:C:252:LEU:C	1:C:254:ILE:H	2.17	0.48
1:C:1:ALA:HB1	1:C:4:GLN:HB3	1.96	0.48
1:B:281:ASN:HB2	1:B:289:ILE:CG2	2.43	0.48
1:B:322:LYS:HG3	1:B:351:PRO:O	2.13	0.48
1:C:20:GLU:CB	1:C:32:ARG:CD	2.92	0.48
1:C:23:ILE:HG23	1:C:29:ASN:ND2	2.28	0.48
1:C:250:TYR:O	1:C:252:LEU:N	2.47	0.48
1:A:354:ALA:HA	1:A:357:ARG:HG3	1.95	0.47
1:C:81:ASN:HD22	1:C:83:HIS:HE1	1.61	0.47
1:B:308:SER:CB	1:B:314:LEU:HD11	2.44	0.47
1:B:13:LEU:HD11	1:B:388:SER:CB	2.43	0.47
1:C:108:PHE:CG	1:C:118:GLY:O	2.67	0.47
1:A:304:LEU:HD12	2:A:506:LDA:H11	1.94	0.47
1:C:139:ALA:HB1	1:C:215:TYR:CE2	2.49	0.47
1:B:9:SER:C	1:B:11:SER:H	2.16	0.47
1:B:159:ALA:O	1:B:160:GLY:O	2.31	0.47
1:A:151:ALA:HB3	1:A:233:ILE:HD13	1.95	0.47
1:C:46:SER:HA	1:C:417:PHE:O	2.14	0.47
1:C:381:ALA:HA	1:C:418:ASN:O	2.13	0.47
1:A:158:PHE:HA	1:A:197:THR:H	1.79	0.47
1:B:10:SER:HB3	1:B:416:ASN:HD22	1.75	0.47
1:B:9:SER:O	1:B:416:ASN:ND2	2.46	0.47
1:C:396:LYS:HD2	1:C:396:LYS:N	2.28	0.47
1:B:124:GLU:HB2	1:B:152:ARG:HB2	1.95	0.47
1:C:51:TYR:HD1	1:C:78:TRP:CD1	2.31	0.47
1:B:169:GLN:HA	1:B:169:GLN:OE1	2.13	0.47
1:C:20:GLU:OE2	1:C:20:GLU:CA	2.60	0.47
1:C:269:LEU:CD2	1:C:271:LEU:HD21	2.45	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:233:ILE:HD12	1:C:269:LEU:HD23	1.97	0.47
1:C:21:GLY:HA2	1:C:36:LEU:CD1	2.35	0.47
1:C:18:SER:OG	1:C:290:HIS:CD2	2.65	0.47
1:A:6:ASN:N	1:A:6:ASN:ND2	2.62	0.47
1:A:284:ASP:CG	1:A:285:PRO:HD2	2.35	0.47
1:A:116:SER:HA	1:A:158:PHE:O	2.14	0.47
1:B:165:LEU:C	1:B:167:ALA:H	2.18	0.47
1:B:334:TYR:N	1:B:334:TYR:CD1	2.83	0.47
1:B:90:ASP:OD2	1:B:90:ASP:N	2.35	0.47
1:A:380:ASP:N	1:A:380:ASP:OD2	2.47	0.47
1:A:304:LEU:CD1	2:A:506:LDA:H11	2.45	0.47
1:B:297:SER:O	1:B:300:GLN:HG3	2.14	0.47
1:A:52:ILE:HG22	1:A:54:PRO:HD3	1.96	0.47
1:A:80:PRO:O	1:A:81:ASN:HB3	2.15	0.47
1:C:317:LYS:HD3	1:C:318:HIS:N	2.30	0.47
1:C:58:ILE:O	1:C:70:ALA:N	2.45	0.47
1:C:64:SER:CB	1:C:165:LEU:HD21	2.34	0.47
1:B:334:TYR:O	1:B:335:TYR:C	2.52	0.47
1:B:216:GLU:HB3	1:B:222:ARG:HB3	1.96	0.47
1:B:235:PHE:HB2	1:B:267:LEU:O	2.15	0.47
1:C:18:SER:HB3	1:C:328:ALA:HB1	1.97	0.46
1:A:304:LEU:O	1:A:316:GLN:HA	2.14	0.46
1:A:256:THR:OG1	1:A:314:LEU:HA	2.15	0.46
1:C:171:MET:HA	1:C:171:MET:HE2	1.96	0.46
1:C:229:SER:C	1:C:230:GLU:O	2.51	0.46
1:A:24:ALA:HB2	1:A:36:LEU:HD22	1.96	0.46
1:C:343:THR:CG2	1:C:344:GLY:N	2.78	0.46
1:A:89:ASN:C	1:A:91:GLN:H	2.19	0.46
1:B:393:GLN:CA	1:B:393:GLN:HE21	2.25	0.46
1:A:374:THR:HG23	1:A:384:ASP:OD2	2.16	0.46
1:C:273:GLU:H	1:C:300:GLN:HE22	1.63	0.46
1:A:162:LEU:HD21	1:A:401:PRO:HG3	1.98	0.46
1:B:354:ALA:HB2	1:B:357:ARG:NH2	2.30	0.46
1:B:191:ILE:HD11	1:B:247:PHE:CZ	2.50	0.46
1:A:1:ALA:HB1	1:A:298:TRP:HE1	1.81	0.46
1:C:26:ASP:O	1:C:28:GLY:N	2.49	0.46
1:A:62:SER:HB3	1:A:68:LEU:HD21	1.96	0.46
1:A:46:SER:HB2	1:A:418:ASN:ND2	2.31	0.46
1:B:137:ASN:N	1:B:137:ASN:HD22	2.13	0.46
1:A:7:GLU:OE2	1:A:52:ILE:HD11	2.16	0.46
1:C:155:ILE:HD11	2:C:501:LDA:C6	2.45	0.46
1:B:354:ALA:HB2	1:B:357:ARG:CZ	2.45	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:286:GLN:O	1:B:333:TYR:HD1	1.99	0.46
1:C:158:PHE:HA	1:C:197:THR:H	1.80	0.46
1:B:19:GLY:N	1:B:292:SER:HB3	2.26	0.46
1:B:191:ILE:HD11	1:B:247:PHE:CE2	2.51	0.46
1:C:317:LYS:HD3	1:C:318:HIS:H	1.80	0.46
1:B:304:LEU:CD1	2:B:503:LDA:C1	2.94	0.46
1:A:62:SER:CB	1:A:68:LEU:HD21	2.46	0.46
1:C:11:SER:HB2	1:C:384:ASP:OD1	2.16	0.45
1:B:228:ARG:HB3	1:B:274:MET:CB	2.46	0.45
1:A:220:ASN:HB3	1:A:282:ARG:HB2	1.99	0.45
1:B:164:GLN:HG3	1:B:189:ASN:ND2	2.31	0.45
1:C:58:ILE:CD1	1:C:406:SER:HB2	2.47	0.45
1:B:155:ILE:HG13	2:B:503:LDA:H92	1.98	0.45
1:C:7:GLU:HB3	1:C:50:VAL:HG21	1.99	0.45
1:B:135:ARG:NH2	1:B:216:GLU:OE2	2.50	0.45
1:C:45:PHE:CG	1:C:46:SER:N	2.84	0.45
1:C:109:ASN:O	1:C:111:THR:N	2.47	0.45
1:C:137:ASN:ND2	1:C:140:TRP:N	2.60	0.45
1:B:162:LEU:O	1:B:166:VAL:HG23	2.16	0.45
1:B:343:THR:HA	1:B:370:SER:O	2.16	0.45
1:A:10:SER:CB	1:A:416:ASN:HD22	2.29	0.45
1:C:51:TYR:CD1	1:C:78:TRP:CD1	3.05	0.45
1:B:114:GLY:HA2	1:B:402:TYR:CZ	2.52	0.45
1:C:123:LEU:HD12	1:C:153:ALA:HB2	1.98	0.45
1:C:361:ILE:CD1	2:C:501:LDA:H22	2.16	0.45
1:A:199:HIS:HE1	1:B:192:ASP:OD1	1.99	0.45
1:A:303:GLN:NE2	1:A:316:GLN:HG3	2.32	0.45
1:A:158:PHE:CA	1:A:197:THR:HB	2.47	0.45
1:C:281:ASN:O	1:C:288:ALA:HA	2.16	0.45
1:B:216:GLU:CB	1:B:222:ARG:HB3	2.47	0.45
1:C:42:ARG:NH1	1:C:420:ALA:HB1	2.31	0.45
1:B:324:ALA:CB	1:B:351:PRO:HG3	2.46	0.45
2:B:503:LDA:H21	2:B:503:LDA:H51	1.29	0.45
1:C:393:GLN:HA	1:C:393:GLN:NE2	2.28	0.45
1:C:334:TYR:CD2	1:C:340:THR:HG23	2.52	0.45
1:A:248:ASN:O	1:A:250:TYR:CD2	2.70	0.45
1:B:378:ASN:O	1:B:379:LYS:CB	2.60	0.45
1:B:121:THR:OG1	1:B:157:ARG:NH2	2.49	0.45
1:A:3:PHE:CD2	1:A:3:PHE:O	2.70	0.45
1:A:68:LEU:HD23	1:A:68:LEU:N	2.32	0.44
1:A:185:ALA:O	1:A:187:THR:N	2.50	0.44
1:C:154:LYS:HA	1:C:200:LEU:O	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:103:GLY:HA2	1:B:123:LEU:O	2.17	0.44
1:A:159:ALA:O	1:A:160:GLY:O	2.35	0.44
1:C:216:GLU:HA	1:C:222:ARG:HB3	1.99	0.44
1:A:5:LEU:HD13	1:A:102:TYR:HA	1.99	0.44
1:B:272:PRO:HG2	1:B:298:TRP:CD2	2.51	0.44
1:A:15:ARG:NH1	1:A:20:GLU:OE1	2.50	0.44
1:C:348:ASP:OD2	1:C:348:ASP:O	2.36	0.44
1:B:129:ASN:CG	1:B:130:LEU:N	2.70	0.44
1:C:343:THR:HG22	1:C:344:GLY:N	2.32	0.44
1:B:305:LYS:HD3	1:B:316:GLN:NE2	2.33	0.44
1:A:51:TYR:OH	1:A:76:THR:HG21	2.18	0.44
1:B:91:GLN:NE2	1:B:135:ARG:O	2.51	0.44
1:C:238:ASN:HD22	1:C:262:THR:CG2	2.30	0.44
1:B:326:ARG:HB2	1:B:348:ASP:CB	2.48	0.44
1:A:361:ILE:O	1:A:361:ILE:HG22	2.17	0.44
1:B:321:PHE:CE2	1:B:352:VAL:HG22	2.53	0.44
1:B:248:ASN:HD21	1:B:256:THR:HA	1.82	0.44
1:C:13:LEU:HD13	1:C:17:TYR:OH	2.18	0.44
1:B:378:ASN:OD1	1:B:380:ASP:N	2.38	0.44
1:B:377:PHE:HB2	1:B:381:ALA:HB3	1.99	0.44
1:A:40:PHE:N	1:A:40:PHE:CD1	2.85	0.44
1:B:254:ILE:HD13	1:B:254:ILE:H	1.83	0.44
1:A:366:ARG:NH1	2:A:505:LDA:H12	2.32	0.44
1:C:27:ALA:O	1:C:28:GLY:C	2.56	0.44
1:A:11:SER:OG	1:A:15:ARG:NH2	2.51	0.44
1:B:343:THR:CG2	1:B:344:GLY:N	2.80	0.44
1:A:216:GLU:HB2	1:A:222:ARG:HB3	2.00	0.44
1:A:15:ARG:NH1	1:A:20:GLU:OE2	2.49	0.44
1:B:280:TYR:CD2	1:B:281:ASN:N	2.86	0.44
1:A:280:TYR:CD2	1:A:280:TYR:C	2.91	0.44
1:C:57:ASN:ND2	1:C:71:ASP:HA	2.33	0.44
1:C:15:ARG:HG2	1:C:342:ARG:HD3	2.00	0.44
1:B:352:VAL:HA	1:B:353:PRO:HD3	1.82	0.44
1:B:304:LEU:HD13	2:B:503:LDA:C1	2.48	0.44
1:A:6:ASN:O	1:A:17:TYR:CE1	2.70	0.44
1:B:165:LEU:C	1:B:167:ALA:N	2.72	0.44
1:A:214:LEU:O	1:A:214:LEU:HD23	2.18	0.44
1:C:1:ALA:HB1	1:C:4:GLN:CB	2.48	0.44
1:B:178:THR:CG2	1:B:181:GLY:H	2.31	0.43
1:B:7:GLU:OE2	2:B:504:LDA:HM12	2.18	0.43
1:B:361:ILE:N	1:B:362:PRO:HD3	2.32	0.43
1:B:155:ILE:HD11	2:B:503:LDA:H71	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:256:THR:OG1	1:B:314:LEU:HA	2.17	0.43
1:B:137:ASN:HD21	1:B:139:ALA:HB3	1.82	0.43
2:C:502:LDA:H12	2:C:502:LDA:H41	1.89	0.43
1:B:113:ALA:HB1	1:B:161:ASP:OD2	2.18	0.43
1:C:181:GLY:C	1:C:183:ALA:H	2.21	0.43
1:B:78:TRP:O	1:B:80:PRO:HD3	2.17	0.43
1:B:36:LEU:HD23	1:B:214:LEU:HD13	2.00	0.43
1:C:195:THR:O	1:C:197:THR:N	2.51	0.43
1:A:36:LEU:O	1:A:39:MET:HB2	2.19	0.43
1:B:13:LEU:HA	1:B:17:TYR:CZ	2.54	0.43
1:A:326:ARG:HH11	1:A:326:ARG:HG3	1.83	0.43
1:A:350:SER:HA	1:A:351:PRO:HD3	1.83	0.43
1:C:345:ILE:CG2	1:C:346:ALA:N	2.82	0.43
1:A:15:ARG:O	1:A:16:ALA:HB3	2.19	0.43
1:B:303:GLN:NE2	1:B:316:GLN:NE2	2.63	0.43
1:C:273:GLU:N	1:C:300:GLN:NE2	2.65	0.43
1:C:334:TYR:CD2	1:C:340:THR:CG2	3.02	0.43
1:A:3:PHE:HD2	1:A:3:PHE:O	2.01	0.43
1:A:34:ALA:HB1	1:A:131:SER:OG	2.18	0.43
1:C:43:PRO:HA	1:C:85:VAL:O	2.18	0.43
1:C:170:ILE:HA	1:C:173:SER:HB3	2.00	0.43
1:A:21:GLY:HA2	1:A:36:LEU:HD11	2.00	0.43
1:C:19:GLY:H	1:C:292:SER:HB3	1.82	0.43
1:B:378:ASN:OD1	1:B:379:LYS:N	2.52	0.43
1:B:13:LEU:HD11	1:B:388:SER:HB2	2.01	0.43
1:C:165:LEU:HD23	1:C:165:LEU:H	1.83	0.43
1:B:13:LEU:CD1	1:B:388:SER:HB2	2.48	0.43
1:C:112:TYR:HE2	1:C:404:PHE:HE2	1.65	0.43
1:C:57:ASN:ND2	1:C:71:ASP:OD2	2.52	0.43
1:A:223:TYR:N	1:A:223:TYR:CD1	2.86	0.43
1:B:252:LEU:HD23	1:B:254:ILE:HG13	2.01	0.43
1:A:5:LEU:CD1	1:A:103:GLY:H	2.32	0.43
1:A:419:TYR:HE2	1:A:421:PHE:CD1	2.37	0.43
1:B:58:ILE:HD13	1:B:406:SER:HB2	1.99	0.43
1:A:86:ALA:HA	1:A:87:PRO:HD3	1.69	0.43
1:C:345:ILE:HG22	1:C:346:ALA:N	2.34	0.43
1:C:321:PHE:CD2	1:C:352:VAL:HG22	2.52	0.43
1:C:404:PHE:CD1	1:C:404:PHE:N	2.86	0.43
1:C:391:HIS:ND1	1:C:392:GLY:N	2.67	0.43
1:C:62:SER:C	1:C:64:SER:N	2.71	0.43
1:C:20:GLU:HB3	1:C:32:ARG:HD2	2.01	0.43
1:C:326:ARG:HG3	1:C:326:ARG:NH1	2.34	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:250:TYR:O	1:B:251:GLY:C	2.57	0.43
1:B:266:TYR:HB2	1:B:307:THR:OG1	2.19	0.42
1:B:10:SER:O	1:B:13:LEU:HB3	2.18	0.42
1:B:55:ASP:O	1:B:408:GLY:HA2	2.19	0.42
1:B:43:PRO:HA	1:B:85:VAL:O	2.19	0.42
1:B:117:VAL:HG23	1:B:359:ILE:HD11	2.01	0.42
1:C:33:ASN:ND2	1:C:226:THR:HG22	2.34	0.42
1:C:243:LEU:HD23	1:C:243:LEU:HA	1.84	0.42
1:B:326:ARG:HA	1:B:348:ASP:CB	2.49	0.42
1:B:290:HIS:HE1	1:B:332:THR:OG1	2.01	0.42
1:B:52:ILE:C	1:B:54:PRO:HD3	2.39	0.42
1:B:280:TYR:C	1:B:280:TYR:HD2	2.22	0.42
1:C:33:ASN:ND2	1:C:226:THR:CG2	2.82	0.42
1:C:20:GLU:HA	1:C:29:ASN:ND2	2.34	0.42
1:C:383:VAL:HG22	1:C:384:ASP:N	2.34	0.42
1:B:165:LEU:O	1:B:167:ALA:N	2.52	0.42
1:C:68:LEU:HD23	1:C:113:ALA:HB3	2.01	0.42
1:C:20:GLU:HB3	1:C:32:ARG:CD	2.50	0.42
1:C:135:ARG:HG3	1:C:137:ASN:O	2.19	0.42
1:A:195:THR:O	1:A:196:LYS:C	2.57	0.42
1:B:58:ILE:HG12	1:B:73:ILE:HD13	2.00	0.42
1:A:220:ASN:HB3	1:A:282:ARG:HB3	2.01	0.42
1:B:249:ASN:O	1:B:250:TYR:CG	2.73	0.42
1:C:8:PHE:C	1:C:8:PHE:CD2	2.92	0.42
1:C:129:ASN:HB2	1:C:147:ASN:OD1	2.19	0.42
1:A:158:PHE:C	1:A:197:THR:HB	2.39	0.42
1:A:33:ASN:C	1:A:35:ALA:N	2.72	0.42
1:C:236:LYS:HD3	1:C:266:TYR:CE2	2.55	0.42
1:C:343:THR:HA	1:C:371:ALA:HA	2.01	0.42
1:C:164:GLN:HG2	1:C:165:LEU:H	1.83	0.42
1:A:168:GLY:O	1:A:169:GLN:C	2.58	0.42
1:B:172:GLN:O	1:B:173:SER:CB	2.68	0.42
1:A:127:ASN:ND2	1:A:149:VAL:HG22	2.34	0.42
1:B:232:LYS:HE2	1:B:270:ASN:ND2	2.34	0.42
1:A:10:SER:HB3	1:A:416:ASN:HD22	1.85	0.42
1:B:137:ASN:N	1:B:137:ASN:ND2	2.67	0.42
1:C:101:ASN:C	1:C:102:TYR:CD1	2.94	0.42
1:B:347:PHE:CE1	1:B:365:ASP:HB3	2.53	0.42
1:B:243:LEU:HD22	1:B:247:PHE:CE1	2.55	0.42
1:A:120:THR:HB	1:A:156:GLU:HB2	2.02	0.42
1:A:153:ALA:HB1	2:A:506:LDA:H82	2.01	0.42
1:B:325:TYR:O	1:B:348:ASP:HB2	2.19	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:121:THR:HG21	2:A:506:LDA:H21	2.02	0.41
1:A:89:ASN:HD22	1:A:90:ASP:N	2.11	0.41
1:C:59:SER:OG	1:C:405:GLU:HB3	2.19	0.41
1:B:32:ARG:HH11	1:B:32:ARG:CG	2.29	0.41
1:B:402:TYR:CB	1:B:404:PHE:CE1	3.03	0.41
1:B:167:ALA:HB1	1:B:185:ALA:O	2.20	0.41
1:B:263:GLN:OE1	1:B:314:LEU:HD21	2.20	0.41
1:C:42:ARG:NH1	1:C:420:ALA:O	2.53	0.41
1:A:404:PHE:CD1	1:A:404:PHE:N	2.89	0.41
1:C:361:ILE:HG22	1:C:361:ILE:O	2.19	0.41
1:A:4:GLN:O	1:A:5:LEU:HD12	2.21	0.41
1:A:15:ARG:HH11	1:A:15:ARG:HG3	1.84	0.41
1:A:30:VAL:HG21	1:A:85:VAL:HG22	2.03	0.41
1:C:117:VAL:HG23	1:C:359:ILE:HD11	2.02	0.41
1:C:218:ASP:C	1:C:218:ASP:OD1	2.59	0.41
1:C:231:VAL:O	1:C:231:VAL:HG12	2.20	0.41
1:C:170:ILE:HG22	1:C:170:ILE:O	2.20	0.41
1:A:6:ASN:HD22	1:A:6:ASN:H	1.69	0.41
1:A:326:ARG:CG	1:A:348:ASP:HB3	2.47	0.41
1:B:342:ARG:NH2	1:B:374:THR:OG1	2.48	0.41
1:A:334:TYR:HA	1:A:340:THR:HG23	2.02	0.41
1:C:133:ALA:HA	1:C:143:GLY:HA2	2.02	0.41
1:C:19:GLY:O	1:C:22:ALA:HB3	2.21	0.41
1:B:112:TYR:OH	1:B:404:PHE:HE2	2.03	0.41
1:A:346:ALA:HB3	1:A:368:TRP:HB2	2.03	0.41
1:A:162:LEU:HD23	1:A:162:LEU:HA	1.90	0.41
1:A:317:LYS:HD3	1:A:318:HIS:N	2.35	0.41
1:B:288:ALA:O	1:B:331:THR:HB	2.21	0.41
1:C:381:ALA:O	1:C:382:SER:HB3	2.21	0.41
1:B:131:SER:HB3	1:B:145:GLY:HA3	2.00	0.41
1:A:19:GLY:O	1:A:20:GLU:C	2.60	0.41
1:B:291:TYR:O	1:B:291:TYR:CD1	2.74	0.41
1:B:321:PHE:N	1:B:321:PHE:CD2	2.89	0.40
1:C:402:TYR:HB2	1:C:404:PHE:CE1	2.55	0.40
1:C:157:ARG:HB2	1:C:197:THR:CG2	2.49	0.40
1:C:366:ARG:NH1	2:C:502:LDA:H21	2.36	0.40
1:B:274:MET:HG2	1:B:296:THR:OG1	2.21	0.40
1:B:347:PHE:HA	1:B:366:ARG:O	2.21	0.40
1:B:128:LEU:HD12	1:B:205:TRP:CZ3	2.56	0.40
1:C:33:ASN:OD1	1:C:33:ASN:C	2.60	0.40
1:A:142:PHE:N	1:A:142:PHE:CD1	2.89	0.40
1:A:356:ASN:O	1:A:357:ARG:C	2.60	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:157:ARG:HH12	2:C:501:LDA:CM2	2.26	0.40
1:B:177:GLN:NE2	1:B:177:GLN:HA	2.35	0.40
1:A:162:LEU:O	1:A:165:LEU:HB2	2.22	0.40
1:A:131:SER:HB2	1:A:144:LEU:O	2.21	0.40
1:A:143:GLY:O	1:A:211:ALA:HA	2.21	0.40
1:B:253:PRO:O	1:B:254:ILE:O	2.40	0.40
1:C:52:ILE:HB	1:C:77:ALA:HB3	2.04	0.40
1:C:140:TRP:O	1:C:141:SER:HB2	2.21	0.40
1:A:158:PHE:N	1:A:158:PHE:CD1	2.89	0.40
1:B:9:SER:C	1:B:11:SER:N	2.75	0.40
1:C:359:ILE:O	1:C:362:PRO:HG3	2.22	0.40
1:B:387:VAL:HA	1:B:412:LEU:O	2.21	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	408/427 (96%)	352 (86%)	43 (10%)	13 (3%)	6	43
1	B	419/427 (98%)	355 (85%)	42 (10%)	22 (5%)	3	26
1	C	412/427 (96%)	340 (82%)	55 (13%)	17 (4%)	4	35
All	All	1239/1281 (97%)	1047 (84%)	140 (11%)	52 (4%)	4	34

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	LYS
1	A	337	ASP
1	A	355	GLN
1	B	90	ASP
1	B	174	PRO
1	B	254	ILE
1	B	355	GLN

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Mol	Chain	Res	Type
1	B	379	LYS
1	C	2	GLY
1	C	27	ALA
1	C	164	GLN
1	C	165	LEU
1	C	355	GLN
1	A	2	GLY
1	A	34	ALA
1	A	160	GLY
1	A	186	ALA
1	B	72	ASN
1	B	160	GLY
1	B	217	LEU
1	B	250	TYR
1	B	272	PRO
1	B	354	ALA
1	B	359	ILE
1	C	29	ASN
1	C	90	ASP
1	C	110	ASP
1	C	160	GLY
1	C	272	PRO
1	A	105	ALA
1	A	249	ASN
1	A	272	PRO
1	B	154	LYS
1	B	362	PRO
1	C	7	GLU
1	C	28	GLY
1	C	196	LYS
1	C	219	LYS
1	C	349	ASP
1	B	365	ASP
1	C	382	SER
1	A	72	ASN
1	B	22	ALA
1	C	335	TYR
1	B	2	GLY
1	B	115	GLY
1	B	335	TYR
1	B	166	VAL
1	B	173	SER

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Mol	Chain	Res	Type
1	B	118	GLY
1	A	63	PRO
1	A	163	GLY

5.3.2 Protein sidechains ⓘ

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	323/336 (96%)	299 (93%)	24 (7%)	20	63
1	B	330/336 (98%)	291 (88%)	39 (12%)	8	34
1	C	326/336 (97%)	293 (90%)	33 (10%)	11	42
All	All	979/1008 (97%)	883 (90%)	96 (10%)	12	45

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	11	SER
1	A	39	MET
1	A	63	PRO
1	A	72	ASN
1	A	78	TRP
1	A	89	ASN
1	A	121	THR
1	A	130	LEU
1	A	137	ASN
1	A	140	TRP
1	A	189	ASN
1	A	197	THR
1	A	203	ASN
1	A	214	LEU
1	A	275	TRP
1	A	278	SER
1	A	282	ARG
1	A	314	LEU
1	A	322	LYS

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Mol	Chain	Res	Type
1	A	380	ASP
1	A	393	GLN
1	A	404	PHE
1	A	416	ASN
1	B	3	PHE
1	B	10	SER
1	B	15	ARG
1	B	32	ARG
1	B	39	MET
1	B	72	ASN
1	B	81	ASN
1	B	89	ASN
1	B	90	ASP
1	B	122	ASP
1	B	125	THR
1	B	137	ASN
1	B	141	SER
1	B	169	GLN
1	B	172	GLN
1	B	174	PRO
1	B	187	THR
1	B	189	ASN
1	B	197	THR
1	B	210	ASN
1	B	249	ASN
1	B	252	LEU
1	B	263	GLN
1	B	264	SER
1	B	275	TRP
1	B	280	TYR
1	B	282	ARG
1	B	290	HIS
1	B	322	LYS
1	B	326	ARG
1	B	334	TYR
1	B	340	THR
1	B	349	ASP
1	B	358	SER
1	B	373	THR
1	B	393	GLN
1	B	403	GLN
1	B	406	SER

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Mol	Chain	Res	Type
1	B	415	THR
1	C	5	LEU
1	C	6	ASN
1	C	39	MET
1	C	63	PRO
1	C	72	ASN
1	C	78	TRP
1	C	120	THR
1	C	121	THR
1	C	122	ASP
1	C	137	ASN
1	C	142	PHE
1	C	144	LEU
1	C	162	LEU
1	C	197	THR
1	C	203	ASN
1	C	214	LEU
1	C	217	LEU
1	C	219	LYS
1	C	263	GLN
1	C	275	TRP
1	C	282	ARG
1	C	284	ASP
1	C	285	PRO
1	C	303	GLN
1	C	322	LYS
1	C	326	ARG
1	C	331	THR
1	C	349	ASP
1	C	355	GLN
1	C	364	GLN
1	C	393	GLN
1	C	396	LYS
1	C	404	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (61) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	29	ASN
1	A	57	ASN
1	A	72	ASN

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Mol	Chain	Res	Type
1	A	89	ASN
1	A	127	ASN
1	A	137	ASN
1	A	147	ASN
1	A	189	ASN
1	A	199	HIS
1	A	210	ASN
1	A	248	ASN
1	A	300	GLN
1	A	303	GLN
1	A	316	GLN
1	A	355	GLN
1	A	356	ASN
1	A	393	GLN
1	A	416	ASN
1	A	418	ASN
1	B	6	ASN
1	B	57	ASN
1	B	72	ASN
1	B	83	HIS
1	B	127	ASN
1	B	129	ASN
1	B	137	ASN
1	B	177	GLN
1	B	189	ASN
1	B	270	ASN
1	B	281	ASN
1	B	290	HIS
1	B	300	GLN
1	B	316	GLN
1	B	356	ASN
1	B	393	GLN
1	B	416	ASN
1	B	418	ASN
1	C	29	ASN
1	C	57	ASN
1	C	72	ASN
1	C	81	ASN
1	C	91	GLN
1	C	129	ASN
1	C	137	ASN
1	C	182	GLN

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Mol	Chain	Res	Type
1	C	189	ASN
1	C	201	ASN
1	C	210	ASN
1	C	238	ASN
1	C	248	ASN
1	C	270	ASN
1	C	281	ASN
1	C	290	HIS
1	C	300	GLN
1	C	316	GLN
1	C	355	GLN
1	C	356	ASN
1	C	393	GLN
1	C	416	ASN
1	C	418	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 ⓘ

6 ligands 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	LDA	A	505	-	15,15,15	3.63	2 (13%)	17,17,17	2.29	6 (35%)
2	LDA	A	506	-	15,15,15	3.82	2 (13%)	17,17,17	2.43	7 (41%)
2	LDA	B	503	-	15,15,15	4.14	2 (13%)	17,17,17	2.25	3 (17%)
2	LDA	B	504	-	15,15,15	3.71	3 (20%)	17,17,17	2.60	5 (29%)
2	LDA	C	501	-	15,15,15	4.07	2 (13%)	17,17,17	2.25	3 (17%)
2	LDA	C	502	-	15,15,15	3.92	2 (13%)	17,17,17	2.53	6 (35%)

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	LDA	A	505	-	-	0/13/13/13	0/0/0/0
2	LDA	A	506	-	-	0/13/13/13	0/0/0/0
2	LDA	B	503	-	-	0/13/13/13	0/0/0/0
2	LDA	B	504	-	-	0/13/13/13	0/0/0/0
2	LDA	C	501	-	-	0/13/13/13	0/0/0/0
2	LDA	C	502	-	-	0/13/13/13	0/0/0/0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	503	LDA	O1-N1	-15.05	1.25	1.39
2	C	501	LDA	O1-N1	-14.84	1.25	1.39
2	A	506	LDA	O1-N1	-13.71	1.26	1.39
2	C	502	LDA	O1-N1	-13.58	1.26	1.39
2	A	505	LDA	O1-N1	-12.66	1.27	1.39
2	B	504	LDA	O1-N1	-12.55	1.27	1.39
2	C	502	LDA	CM2-N1	-5.76	1.40	1.49
2	B	504	LDA	CM2-N1	-5.43	1.41	1.49
2	A	506	LDA	CM2-N1	-5.10	1.41	1.49
2	A	505	LDA	CM2-N1	-5.07	1.41	1.49
2	C	501	LDA	CM2-N1	-4.66	1.42	1.49
2	B	503	LDA	CM2-N1	-4.66	1.42	1.49
2	B	504	LDA	CM1-N1	-3.10	1.44	1.49

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	504	LDA	CM2-N1-CM1	-7.95	99.77	108.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	506	LDA	CM2-N1-CM1	-7.04	100.80	108.85
2	C	502	LDA	CM2-N1-CM1	-6.93	100.93	108.85
2	B	503	LDA	CM2-N1-CM1	-6.88	100.98	108.85
2	C	501	LDA	CM2-N1-CM1	-6.43	101.50	108.85
2	A	505	LDA	CM2-N1-CM1	-6.25	101.70	108.85
2	C	502	LDA	O1-N1-C1	4.62	116.62	110.19
2	C	501	LDA	C2-C1-N1	4.55	121.64	113.80
2	B	504	LDA	O1-N1-C1	4.17	116.00	110.19
2	A	505	LDA	O1-N1-C1	4.09	115.89	110.19
2	B	503	LDA	O1-N1-CM2	3.84	114.18	109.01
2	A	506	LDA	O1-N1-C1	3.32	114.82	110.19
2	C	502	LDA	O1-N1-CM2	3.26	113.40	109.01
2	B	504	LDA	O1-N1-CM2	3.13	113.23	109.01
2	A	506	LDA	O1-N1-CM2	3.08	113.16	109.01
2	B	504	LDA	CM1-N1-C1	-2.97	100.13	109.86
2	A	505	LDA	O1-N1-CM2	2.94	112.97	109.01
2	B	503	LDA	O1-N1-C1	2.78	114.06	110.19
2	C	502	LDA	O1-N1-CM1	-2.66	105.43	109.01
2	A	505	LDA	CM1-N1-C1	-2.46	101.82	109.86
2	C	502	LDA	C2-C1-N1	2.43	117.99	113.80
2	A	506	LDA	CM1-N1-C1	-2.41	101.98	109.86
2	C	501	LDA	O1-N1-CM2	2.35	112.17	109.01
2	A	505	LDA	C2-C1-N1	2.29	117.75	113.80
2	A	506	LDA	C6-C5-C4	-2.21	102.64	114.61
2	C	502	LDA	C9-C8-C7	-2.20	102.72	114.61
2	A	505	LDA	C9-C8-C7	-2.14	103.02	114.61
2	A	506	LDA	C9-C8-C7	-2.14	103.05	114.61
2	B	504	LDA	C2-C1-N1	2.13	117.47	113.80
2	A	506	LDA	C2-C1-N1	2.01	117.27	113.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	412/427 (96%)	-0.20	1 (0%) 93 70	4, 20, 35, 50	0
1	B	421/427 (98%)	-0.18	0 100 100	10, 24, 35, 46	0
1	C	416/427 (97%)	-0.17	0 100 100	7, 25, 38, 45	0
All	All	1249/1281 (97%)	-0.19	1 (0%) 93 77	4, 23, 36, 50	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	184	LEU	2.2

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	LDA	B	504	16/16	0.36	5.49	22,32,37,37	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	LDA	C	501	16/16	0.39	3.55	18,21,31,31	0
2	LDA	C	502	16/16	0.56	3.42	30,37,47,47	0
2	LDA	A	505	16/16	0.37	3.36	25,29,30,30	0
2	LDA	A	506	16/16	0.36	3.06	21,30,46,46	0
2	LDA	B	503	16/16	0.32	2.35	23,26,39,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.