



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 01:35 AM GMT

PDB ID : 1S1U
Title : Crystal structure of L100I mutant HIV-1 reverse transcriptase in complex with nevirapine
Authors : Ren, J.; Nichols, C.E.; Chamberlain, P.P.; Stammers, D.K.
Deposited on : 2004-01-07
Resolution : 3.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

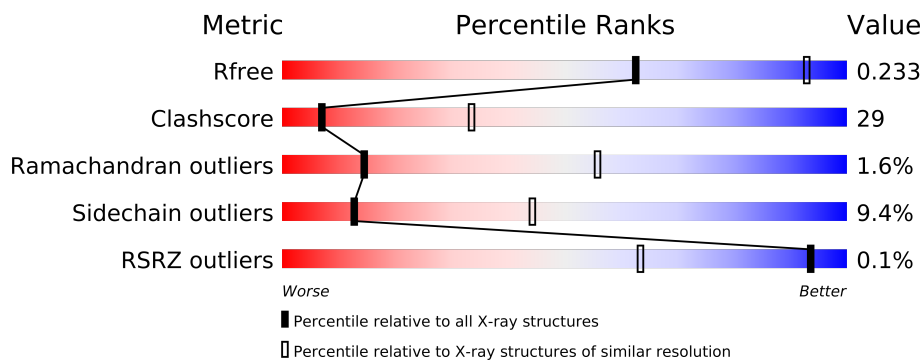
The following versions of software and data (see [references](#)) were used in the production of this report:

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

1 Overall quality at a glance

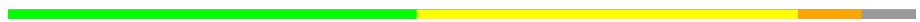
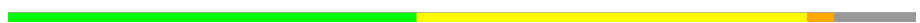
The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	560	
2	B	440	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7660 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 Reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	525	Total	C	N	O	S	0	0	0
			4310	2792	714	796	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ILE	LEU	ENGINEERED	UNP P04585
A	280	CSD	CYS	MODIFIED RESIDUE	UNP P04585

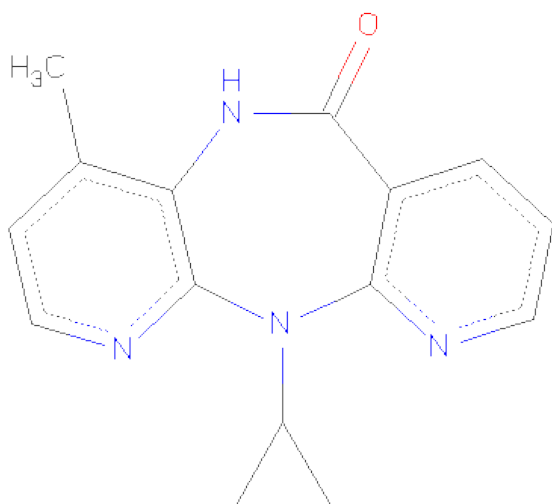
- Molecule 2 is a protein called Reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	401	Total	C	N	O	S	0	0	0
			3330	2173	549	601	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	100	ILE	LEU	ENGINEERED	UNP P04585

- Molecule 3 is 11-CYCLOPROPYL-5,11-DIHYDRO-4-METHYL-6H-DIPYRIDO[3,2-B:2',3'-E][1,4]DIAZEPIN-6-ONE (three-letter code: NVP) (formula: C₁₅H₁₄N₄O).



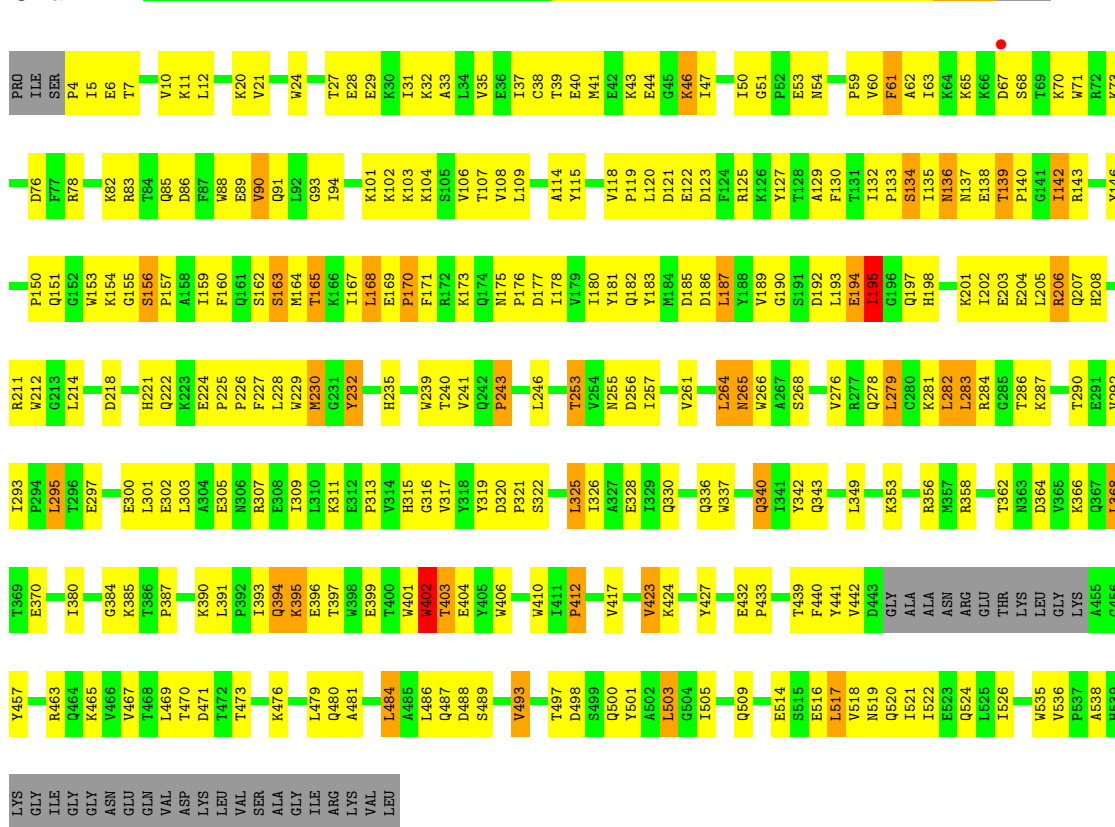
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	20	15	4	1	0	0

3 Residue-property plots

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

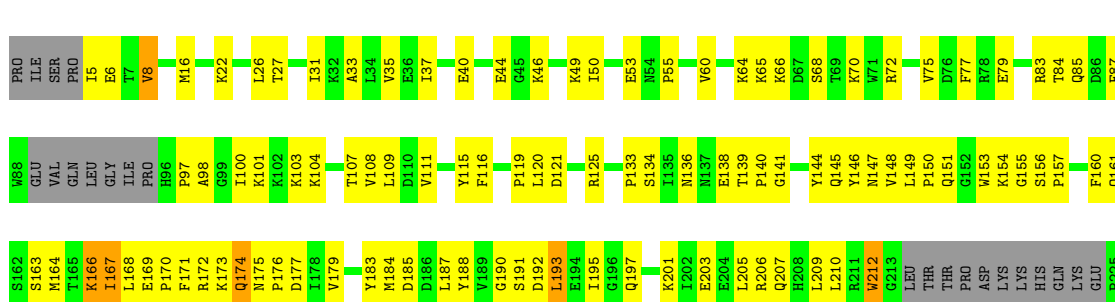
• Molecule 1: Reverse transcriptase

Chain A:



• Molecule 2: Reverse transcriptase

Chain B:



G384	K385	T386	P387	K388	F389	K390	T393	Q394	K395	K398	W398	W401	W402	T403	E404	Y405	W406	Q407	W414	E415	F416	W417	N418	T419	P420	P421	L422	V423	K424	L425	W426	Y427	Q428	LEU	GLU	LYS	GLU	PRO	ILE	VAL	GLY	ALA	GLU	THR	PHE								
P313	G316	V317	Y318	K323	A327	E328	I329	Q330	F331	Q332	G335	Q336	W337	T338	Y339	Q340	I341	Y342	Q343	E344	P345	F346	K347	N348	L349	K350	T351	G352	K353	Y354	A355	R356	MET	ARG	GLY	ALA	HIS	T362	N363	D364	V365	K366	Q367	L368	T369	E370	A371	V372	I375	S379	I380	K311	W383
P226	F227	L228	W229	W230	Q231	Y232	E233	L234	W239	T240	Y241	P247	E248	K249	W252	T253	W254	Q258	V261	G262	K263	L264	W265	W266	W267	S268	Q269	L270	Y276	R277	C280	L283	T286	L295	T296	E297	E298	L301	E302	L303	A304	E305	N306	I309	L310	K311	E312						

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	139.80Å 115.50Å 66.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.90 – 3.00 29.90 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.5 (29.90-3.00) 98.6 (29.90-3.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 3.00Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.227 , 0.285 0.216 , 0.233	Depositor DCC
R_{free} test set	1060 reflections (5.13%)	DCC
Wilson B-factor (Å ²)	84.0	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 76.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 21742 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7660	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: CSD, NVP

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	0.52	0/4417	0.75	1/6005 (0.0%)
2	B	0.51	0/3426	0.73	1/4653 (0.0%)
All	All	0.52	0/7843	0.74	2/10658 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	VAL	N-CA-C	-5.53	96.06	111.00
2	B	229	TRP	CA-CB-CG	-5.09	104.02	113.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4310	0	4343	243	0
2	B	3330	0	3353	207	0
3	A	20	0	14	0	0
All	All	7660	0	7710	440	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 29.

All (440) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:469:LEU:HD21	1:A:480:GLN:HG3	1.17	1.10
2:B:66:LYS:HG2	2:B:230:MET:HA	1.35	1.07
1:A:195:ILE:H	1:A:195:ILE:HD13	1.25	1.01
1:A:295:LEU:HB3	1:A:300:GLU:HG3	1.46	0.96
1:A:517:LEU:HA	1:A:520:GLN:HE21	1.30	0.95
2:B:227:PHE:HB2	2:B:231:GLY:HA2	1.49	0.94
2:B:167:ILE:HD12	2:B:212:TRP:HB3	1.46	0.94
1:A:228:LEU:HD23	1:A:228:LEU:H	1.32	0.93
2:B:227:PHE:CB	2:B:231:GLY:HA2	2.00	0.91
2:B:166:LYS:HA	2:B:166:LYS:HE3	1.51	0.91
1:A:167:ILE:O	1:A:170:PRO:HD2	1.73	0.88
1:A:469:LEU:HD21	1:A:480:GLN:CG	2.04	0.87
1:A:171:PHE:CE1	1:A:205:LEU:HA	2.11	0.85
2:B:297:GLU:O	2:B:301:LEU:HD22	1.76	0.85
1:A:399:GLU:HA	1:A:402:TRP:HE3	1.42	0.84
1:A:46:LYS:HB3	1:A:46:LYS:NZ	1.93	0.84
2:B:66:LYS:CG	2:B:230:MET:HA	2.08	0.83
2:B:66:LYS:HE3	2:B:230:MET:CG	2.08	0.83
1:A:65:LYS:HB3	1:A:68:SER:HB3	1.59	0.82
1:A:226:PRO:HG3	1:A:235:HIS:CE1	2.16	0.80
1:A:142:ILE:H	1:A:142:ILE:HD13	1.46	0.80
2:B:66:LYS:HG2	2:B:230:MET:CA	2.10	0.80
2:B:230:MET:C	2:B:232:TYR:H	1.81	0.79
2:B:227:PHE:HB2	2:B:231:GLY:CA	2.12	0.79
1:A:469:LEU:HD11	1:A:480:GLN:HE21	1.48	0.79
1:A:165:THR:HG21	2:B:140:PRO:HG2	1.65	0.79
1:A:469:LEU:CD2	1:A:480:GLN:HG3	2.07	0.78
2:B:175:ASN:ND2	2:B:201:LYS:NZ	2.33	0.77
2:B:173:LYS:HA	2:B:176:PRO:HG3	1.66	0.77
1:A:169:GLU:HB3	1:A:170:PRO:HD3	1.66	0.76
2:B:175:ASN:HD21	2:B:201:LYS:HZ1	1.31	0.76
1:A:206:ARG:NH1	1:A:218:ASP:HA	2.00	0.75
2:B:167:ILE:HD12	2:B:212:TRP:CB	2.16	0.75
2:B:175:ASN:HD21	2:B:201:LYS:NZ	1.84	0.75
1:A:156:SER:HB2	1:A:157:PRO:HD3	1.67	0.75
1:A:328:GLU:HG3	1:A:390:LYS:HB2	1.68	0.75
2:B:111:VAL:HG11	2:B:187:LEU:HD22	1.69	0.75
2:B:103:LYS:NZ	2:B:191:SER:HA	2.02	0.74
2:B:169:GLU:N	2:B:170:PRO:HD2	2.02	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:27:THR:HG22	1:A:29:GLU:H	1.53	0.73
2:B:295:LEU:HD12	2:B:295:LEU:H	1.53	0.73
2:B:395:LYS:HD2	2:B:416:PHE:CE1	2.23	0.73
2:B:420:PRO:O	2:B:423:VAL:HG12	1.88	0.73
1:A:325:LEU:HD22	1:A:385:LYS:HE3	1.70	0.73
2:B:353:LYS:NZ	2:B:428:GLN:HG3	2.03	0.73
2:B:332:GLN:HB3	2:B:428:GLN:HE22	1.54	0.72
2:B:136:ASN:HB3	2:B:138:GLU:HG3	1.71	0.72
1:A:406:TRP:CH2	2:B:418:ASN:HA	2.24	0.72
1:A:101:LYS:H	1:A:101:LYS:HD2	1.55	0.71
2:B:254:VAL:O	2:B:258:GLN:HG3	1.92	0.70
2:B:229:TRP:CE3	2:B:229:TRP:HA	2.26	0.70
1:A:134:SER:CB	1:A:139:THR:O	2.40	0.69
2:B:154:LYS:O	2:B:157:PRO:HD2	1.92	0.69
1:A:228:LEU:CD2	1:A:228:LEU:H	2.04	0.69
1:A:295:LEU:HB3	1:A:300:GLU:CG	2.20	0.69
2:B:261:VAL:HG13	2:B:276:VAL:HG11	1.75	0.69
1:A:253:THR:HG22	1:A:256:ASP:H	1.55	0.69
2:B:295:LEU:HD12	2:B:295:LEU:N	2.08	0.69
1:A:101:LYS:N	1:A:101:LYS:HD2	2.07	0.69
1:A:325:LEU:CD2	1:A:385:LYS:HE3	2.22	0.68
1:A:295:LEU:HD12	1:A:300:GLU:HG2	1.75	0.68
2:B:139:THR:HG23	2:B:140:PRO:HD2	1.75	0.68
1:A:21:VAL:HG13	1:A:59:PRO:HD3	1.76	0.67
2:B:175:ASN:CG	2:B:201:LYS:HZ2	1.96	0.67
1:A:134:SER:OG	1:A:140:PRO:HA	1.92	0.67
2:B:266:TRP:O	2:B:269:GLN:HG2	1.94	0.67
1:A:165:THR:HG23	1:A:182:GLN:OE1	1.94	0.67
1:A:194:GLU:OE1	1:A:197:GLN:HB2	1.95	0.67
2:B:66:LYS:HE3	2:B:230:MET:HA	1.77	0.66
1:A:295:LEU:CB	1:A:300:GLU:HG3	2.23	0.66
1:A:115:TYR:OH	1:A:157:PRO:HG3	1.94	0.66
1:A:228:LEU:HD23	1:A:228:LEU:N	2.09	0.66
1:A:12:LEU:HD11	1:A:127:TYR:CE1	2.31	0.66
1:A:180:ILE:HG12	1:A:189:VAL:HG13	1.76	0.66
1:A:142:ILE:H	1:A:142:ILE:CD1	1.99	0.66
1:A:399:GLU:HA	1:A:402:TRP:CE3	2.30	0.66
1:A:457:TYR:HE1	1:A:463:ARG:HG2	1.60	0.66
2:B:156:SER:HB2	2:B:157:PRO:HD3	1.78	0.65
2:B:97:PRO:O	2:B:100:ILE:HG22	1.96	0.65
2:B:31:ILE:CD1	2:B:133:PRO:HG2	2.26	0.65
2:B:393:ILE:HG12	2:B:394:GLN:N	2.11	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:227:PHE:CB	2:B:231:GLY:CA	2.73	0.65
1:A:132:ILE:HB	1:A:142:ILE:HG12	1.78	0.64
2:B:103:LYS:HZ1	2:B:191:SER:HA	1.62	0.64
2:B:65:LYS:HB2	2:B:68:SER:HB3	1.79	0.64
2:B:66:LYS:HE3	2:B:230:MET:CB	2.27	0.64
1:A:358:ARG:NH2	2:B:394:GLN:OE1	2.29	0.64
2:B:372:VAL:HA	2:B:389:PHE:CE2	2.33	0.64
2:B:5:ILE:HG22	2:B:6:GLU:N	2.11	0.64
2:B:372:VAL:HA	2:B:389:PHE:HE2	1.63	0.63
1:A:395:LYS:CD	1:A:395:LYS:H	2.10	0.63
1:A:181:TYR:CE1	1:A:183:TYR:HB2	2.34	0.63
1:A:83:ARG:HG3	1:A:83:ARG:HH11	1.64	0.63
1:A:171:PHE:HE1	1:A:205:LEU:HA	1.65	0.62
1:A:232:TYR:HD2	1:A:239:TRP:HZ3	1.47	0.62
2:B:183:TYR:CE1	2:B:184:MET:HG2	2.34	0.62
1:A:201:LYS:HA	1:A:204:GLU:HG3	1.81	0.62
1:A:46:LYS:HB3	1:A:46:LYS:HZ2	1.62	0.62
2:B:66:LYS:CE	2:B:230:MET:HA	2.30	0.61
2:B:393:ILE:HG12	2:B:394:GLN:H	1.65	0.61
1:A:195:ILE:HD13	1:A:195:ILE:N	2.07	0.61
1:A:279:LEU:HD22	1:A:302:GLU:OE1	2.00	0.61
1:A:498:ASP:HA	1:A:536:VAL:O	1.99	0.61
2:B:301:LEU:HD13	2:B:301:LEU:N	2.16	0.61
2:B:203:GLU:O	2:B:206:ARG:HB2	2.00	0.61
2:B:5:ILE:HG22	2:B:6:GLU:H	1.66	0.60
1:A:104:LYS:HB2	1:A:192:ASP:HA	1.83	0.60
2:B:134:SER:OG	2:B:139:THR:HB	2.01	0.60
2:B:348:ASN:HD22	2:B:351:THR:CG2	2.15	0.60
1:A:524:GLN:HA	1:A:524:GLN:OE1	2.00	0.60
1:A:393:ILE:HB	1:A:423:VAL:HG22	1.84	0.60
1:A:465:LYS:HD3	1:A:488:ASP:OD1	2.01	0.59
1:A:330:GLN:NE2	1:A:340:GLN:OE1	2.30	0.59
2:B:193:LEU:HD23	2:B:197:GLN:HB3	1.82	0.59
2:B:100:ILE:HG23	2:B:101:LYS:N	2.17	0.59
1:A:279:LEU:HA	1:A:282:LEU:CD2	2.32	0.59
1:A:171:PHE:CE1	1:A:205:LEU:CA	2.83	0.59
2:B:183:TYR:CD1	2:B:184:MET:HG2	2.37	0.59
2:B:366:LYS:HD2	2:B:405:TYR:CE2	2.37	0.59
1:A:218:ASP:OD1	1:A:221:HIS:CD2	2.55	0.59
1:A:325:LEU:H	1:A:325:LEU:HD23	1.67	0.59
1:A:239:TRP:NE1	1:A:316:GLY:HA3	2.17	0.59
2:B:379:SER:OG	2:B:387:PRO:HG3	2.02	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:295:LEU:CB	1:A:300:GLU:CG	2.80	0.58
2:B:340:GLN:HB3	2:B:348:ASN:ND2	2.18	0.58
1:A:61:PHE:N	1:A:61:PHE:CD2	2.72	0.58
1:A:136:ASN:HB3	1:A:139:THR:OG1	2.04	0.58
2:B:66:LYS:HE3	2:B:230:MET:HG3	1.84	0.58
1:A:253:THR:HG23	1:A:255:ASN:H	1.68	0.58
1:A:516:GLU:O	1:A:520:GLN:HG3	2.03	0.58
1:A:342:TYR:HA	1:A:349:LEU:HD23	1.84	0.58
1:A:193:LEU:HD12	1:A:198:HIS:HA	1.86	0.58
1:A:195:ILE:H	1:A:195:ILE:CD1	1.93	0.58
2:B:203:GLU:HG3	2:B:207:GLN:HE22	1.69	0.57
2:B:175:ASN:CG	2:B:201:LYS:NZ	2.58	0.57
2:B:227:PHE:HB3	2:B:231:GLY:HA2	1.85	0.57
2:B:163:SER:O	2:B:167:ILE:HG22	2.04	0.57
1:A:311:LYS:O	1:A:313:PRO:HD3	2.05	0.57
2:B:305:GLU:O	2:B:309:ILE:HG13	2.04	0.57
1:A:412:PRO:HG3	2:B:401:TRP:HZ2	1.69	0.57
2:B:234:LEU:HD23	2:B:239:TRP:CZ2	2.39	0.57
2:B:139:THR:HG23	2:B:140:PRO:CD	2.35	0.57
1:A:229:TRP:O	1:A:232:TYR:HD1	1.87	0.56
2:B:175:ASN:OD1	2:B:201:LYS:NZ	2.36	0.56
1:A:91:GLN:NE2	2:B:140:PRO:O	2.37	0.56
2:B:328:GLU:HG2	2:B:390:LYS:HD2	1.87	0.56
1:A:106:VAL:HG22	1:A:190:GLY:HA3	1.87	0.56
1:A:10:VAL:HG12	1:A:11:LYS:N	2.21	0.56
1:A:305:GLU:O	1:A:309:ILE:HG13	2.05	0.56
1:A:115:TYR:HD1	1:A:151:GLN:HA	1.71	0.56
1:A:358:ARG:CZ	2:B:394:GLN:OE1	2.54	0.56
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.86	0.56
2:B:241:VAL:HG11	2:B:313:PRO:HG3	1.88	0.56
1:A:169:GLU:O	1:A:173:LYS:HG3	2.05	0.56
2:B:139:THR:HG22	2:B:141:GLY:H	1.70	0.56
1:A:257:ILE:O	1:A:261:VAL:HG23	2.06	0.55
1:A:390:LYS:O	1:A:391:LEU:HD23	2.06	0.55
2:B:50:ILE:CG2	2:B:145:GLN:HG2	2.36	0.55
1:A:5:ILE:CG1	1:A:6:GLU:N	2.69	0.55
2:B:295:LEU:H	2:B:295:LEU:CD1	2.19	0.55
1:A:61:PHE:N	1:A:61:PHE:HD2	2.04	0.55
2:B:371:ALA:O	2:B:375:ILE:HG13	2.06	0.55
2:B:169:GLU:N	2:B:170:PRO:CD	2.70	0.55
2:B:372:VAL:HG13	2:B:389:PHE:CE2	2.42	0.55
2:B:369:THR:HG22	2:B:398:TRP:CH2	2.42	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:106:VAL:HA	1:A:190:GLY:HA2	1.87	0.55
1:A:320:ASP:O	1:A:343:GLN:NE2	2.37	0.55
1:A:228:LEU:HG	1:A:228:LEU:O	2.07	0.54
1:A:229:TRP:O	1:A:230:MET:C	2.45	0.54
1:A:229:TRP:O	1:A:232:TYR:CD1	2.60	0.54
1:A:138:GLU:HG3	1:A:138:GLU:O	2.07	0.54
1:A:39:THR:O	1:A:43:LYS:HG2	2.06	0.54
1:A:412:PRO:HG3	2:B:401:TRP:CZ2	2.41	0.54
1:A:264:LEU:HD23	1:A:276:VAL:HG12	1.90	0.54
1:A:115:TYR:HE2	1:A:160:PHE:CD1	2.26	0.54
1:A:37:ILE:O	1:A:41:MET:HG3	2.08	0.54
2:B:65:LYS:CB	2:B:68:SER:HB3	2.37	0.54
2:B:230:MET:HB2	2:B:232:TYR:HB2	1.89	0.54
2:B:372:VAL:HG13	2:B:389:PHE:CZ	2.42	0.54
1:A:239:TRP:O	1:A:315:HIS:HA	2.07	0.54
1:A:27:THR:O	1:A:31:ILE:HG13	2.08	0.54
2:B:332:GLN:HB2	2:B:336:GLN:O	2.08	0.54
2:B:66:LYS:HE3	2:B:230:MET:CA	2.38	0.53
2:B:79:GLU:O	2:B:83:ARG:HG3	2.08	0.53
1:A:159:ILE:O	1:A:162:SER:HB3	2.09	0.53
1:A:268:SER:HB3	1:A:353:LYS:HE2	1.89	0.53
2:B:173:LYS:O	2:B:176:PRO:HD3	2.07	0.53
1:A:134:SER:HB3	1:A:139:THR:O	2.07	0.53
1:A:229:TRP:N	1:A:232:TYR:O	2.41	0.53
2:B:49:LYS:HG3	2:B:144:TYR:CE1	2.43	0.53
1:A:165:THR:HA	1:A:168:LEU:HD22	1.90	0.53
1:A:78:ARG:O	1:A:82:LYS:HG3	2.09	0.53
1:A:395:LYS:HD2	1:A:395:LYS:N	2.24	0.53
1:A:467:VAL:HG21	1:A:484:LEU:HD11	1.90	0.53
2:B:175:ASN:N	2:B:176:PRO:HD3	2.24	0.52
1:A:297:GLU:O	1:A:301:LEU:HB2	2.08	0.52
1:A:146:TYR:CG	1:A:150:PRO:HG3	2.44	0.52
1:A:38:CYS:O	1:A:47:ILE:HD11	2.10	0.52
1:A:417:VAL:O	1:A:417:VAL:HG13	2.09	0.52
2:B:385:LYS:HG2	2:B:386:THR:N	2.24	0.52
2:B:139:THR:HG22	2:B:141:GLY:N	2.25	0.52
2:B:366:LYS:O	2:B:370:GLU:HG3	2.10	0.51
1:A:11:LYS:O	1:A:85:GLN:HG2	2.10	0.51
2:B:302:GLU:O	2:B:306:ASN:ND2	2.43	0.51
1:A:31:ILE:HG12	1:A:133:PRO:HG2	1.91	0.51
1:A:319:TYR:CD2	1:A:320:ASP:N	2.78	0.51
2:B:249:LYS:HG3	2:B:252:TRP:CE2	2.45	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:31:ILE:HD11	2:B:133:PRO:HG2	1.91	0.51
1:A:63:ILE:N	1:A:63:ILE:HD12	2.25	0.51
1:A:5:ILE:HG12	1:A:6:GLU:N	2.25	0.51
1:A:142:ILE:N	1:A:142:ILE:HD13	2.20	0.51
2:B:174:GLN:NE2	2:B:175:ASN:OD1	2.44	0.51
1:A:94:ILE:HG23	1:A:229:TRP:CH2	2.45	0.51
1:A:503:LEU:HG	1:A:535:TRP:HB2	1.91	0.51
1:A:46:LYS:HB3	1:A:46:LYS:HZ3	1.72	0.51
1:A:24:TRP:HZ3	1:A:61:PHE:CD2	2.28	0.51
2:B:161:GLN:O	2:B:164:MET:HB3	2.10	0.51
2:B:139:THR:CG2	2:B:140:PRO:N	2.73	0.51
1:A:76:ASP:OD2	1:A:78:ARG:HG3	2.11	0.51
2:B:423:VAL:HG13	2:B:424:LYS:N	2.26	0.51
2:B:342:TYR:HB3	2:B:348:ASN:HA	1.92	0.51
2:B:40:GLU:O	2:B:44:GLU:HG3	2.10	0.51
2:B:206:ARG:HD2	2:B:227:PHE:CE1	2.45	0.51
1:A:240:THR:OG1	1:A:241:VAL:N	2.44	0.51
1:A:86:ASP:HA	1:A:154:LYS:HE2	1.93	0.51
2:B:103:LYS:HE3	2:B:190:GLY:C	2.32	0.50
1:A:457:TYR:CE1	1:A:463:ARG:HG2	2.44	0.50
1:A:326:ILE:HB	1:A:342:TYR:CE1	2.46	0.50
2:B:380:ILE:O	2:B:384:GLY:HA2	2.10	0.50
2:B:227:PHE:HB3	2:B:231:GLY:O	2.12	0.50
1:A:395:LYS:HD2	1:A:395:LYS:H	1.77	0.50
1:A:108:VAL:HG23	1:A:227:PHE:CZ	2.47	0.50
2:B:402:TRP:CE2	2:B:403:THR:HG23	2.47	0.50
2:B:103:LYS:HE2	2:B:179:VAL:HG23	1.92	0.50
2:B:350:LYS:HG3	2:B:351:THR:N	2.27	0.50
1:A:232:TYR:HD2	1:A:239:TRP:CZ3	2.29	0.50
2:B:85:GLN:O	2:B:87:PHE:CE1	2.65	0.50
2:B:195:ILE:HD11	2:B:233:GLU:CD	2.32	0.50
1:A:337:TRP:CZ3	1:A:368:LEU:HD23	2.47	0.50
2:B:206:ARG:O	2:B:210:LEU:HD23	2.11	0.50
1:A:481:ALA:O	1:A:484:LEU:HB2	2.11	0.50
1:A:122:GLU:HA	1:A:125:ARG:HD2	1.94	0.49
2:B:298:GLU:OE1	2:B:298:GLU:N	2.44	0.49
2:B:134:SER:CB	2:B:139:THR:HB	2.42	0.49
2:B:330:GLN:HB2	2:B:338:THR:OG1	2.12	0.49
2:B:64:LYS:HE2	2:B:68:SER:O	2.12	0.49
1:A:358:ARG:NE	2:B:394:GLN:HE22	2.11	0.49
2:B:311:LYS:O	2:B:312:GLU:HG3	2.13	0.49
2:B:301:LEU:N	2:B:301:LEU:CD1	2.76	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:175:ASN:ND2	2:B:201:LYS:CE	2.75	0.49
2:B:353:LYS:HZ2	2:B:428:GLN:HG3	1.76	0.49
1:A:181:TYR:HE1	1:A:183:TYR:HB2	1.78	0.49
1:A:402:TRP:CG	1:A:403:THR:N	2.81	0.49
1:A:24:TRP:HZ3	1:A:61:PHE:CG	2.30	0.49
2:B:195:ILE:HD11	2:B:233:GLU:HG2	1.95	0.49
1:A:295:LEU:HB2	1:A:300:GLU:OE1	2.13	0.48
1:A:497:THR:O	1:A:535:TRP:HA	2.13	0.48
1:A:186:ASP:O	1:A:187:LEU:HD22	2.13	0.48
1:A:62:ALA:C	1:A:63:ILE:HD12	2.33	0.48
2:B:44:GLU:HB2	2:B:46:LYS:HG2	1.95	0.48
1:A:50:ILE:HD12	1:A:54:ASN:HB3	1.96	0.48
2:B:423:VAL:CG1	2:B:424:LYS:N	2.75	0.48
2:B:100:ILE:CG2	2:B:101:LYS:N	2.75	0.48
1:A:432:GLU:HB2	1:A:433:PRO:HD2	1.95	0.48
1:A:469:LEU:HD11	1:A:480:GLN:NE2	2.23	0.48
1:A:106:VAL:HG12	1:A:107:THR:N	2.29	0.48
1:A:380:ILE:O	1:A:384:GLY:HA2	2.13	0.48
1:A:107:THR:HG22	1:A:109:LEU:CD1	2.43	0.48
2:B:425:LEU:HD22	2:B:426:TRP:CD1	2.49	0.48
2:B:167:ILE:CG1	2:B:167:ILE:O	2.61	0.47
2:B:191:SER:HB2	2:B:193:LEU:HD13	1.96	0.47
1:A:202:ILE:HG22	1:A:203:GLU:N	2.29	0.47
1:A:175:ASN:N	1:A:176:PRO:HD3	2.28	0.47
1:A:279:LEU:O	1:A:282:LEU:HB2	2.14	0.47
2:B:380:ILE:O	2:B:384:GLY:N	2.45	0.47
2:B:175:ASN:ND2	2:B:201:LYS:HZ1	2.00	0.47
1:A:279:LEU:HA	1:A:282:LEU:HD22	1.94	0.47
2:B:87:PHE:CD1	2:B:87:PHE:N	2.81	0.47
1:A:206:ARG:CZ	1:A:218:ASP:HA	2.44	0.47
1:A:203:GLU:O	1:A:204:GLU:C	2.53	0.47
1:A:155:GLY:O	1:A:156:SER:C	2.53	0.47
1:A:101:LYS:N	1:A:101:LYS:CD	2.77	0.47
1:A:232:TYR:CD2	1:A:239:TRP:CZ3	3.03	0.47
1:A:28:GLU:O	1:A:32:LYS:HG3	2.12	0.47
1:A:168:LEU:N	1:A:168:LEU:HD13	2.29	0.47
2:B:379:SER:HA	2:B:383:TRP:CE3	2.49	0.47
2:B:234:LEU:HD23	2:B:239:TRP:CH2	2.50	0.47
1:A:320:ASP:OD2	1:A:322:SER:OG	2.28	0.47
2:B:422:LEU:O	2:B:425:LEU:HB3	2.14	0.47
1:A:303:LEU:HD23	1:A:303:LEU:O	2.14	0.47
2:B:167:ILE:HD11	2:B:209:LEU:HD23	1.97	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:325:LEU:CD2	1:A:325:LEU:H	2.28	0.47
1:A:28:GLU:HG3	1:A:135:ILE:HG21	1.96	0.47
2:B:366:LYS:HD2	2:B:405:TYR:CD2	2.49	0.47
1:A:427:TYR:OH	1:A:509:GLN:HA	2.15	0.47
1:A:522:ILE:O	1:A:526:ILE:HG13	2.14	0.47
1:A:265:ASN:O	1:A:266:TRP:C	2.52	0.47
2:B:206:ARG:HD2	2:B:227:PHE:HE1	1.80	0.46
2:B:66:LYS:CD	2:B:230:MET:HA	2.45	0.46
2:B:365:VAL:O	2:B:366:LYS:C	2.52	0.46
2:B:66:LYS:HE3	2:B:230:MET:HG2	1.96	0.46
2:B:332:GLN:HB3	2:B:428:GLN:NE2	2.27	0.46
2:B:345:PRO:C	2:B:347:LYS:H	2.18	0.46
2:B:116:PHE:CZ	2:B:151:GLN:HG3	2.51	0.46
1:A:208:HIS:O	1:A:212:TRP:CD1	2.69	0.46
1:A:486:LEU:CD1	1:A:521:ILE:HG23	2.45	0.46
2:B:175:ASN:ND2	2:B:201:LYS:HE3	2.31	0.46
1:A:439:THR:HG22	1:A:441:TYR:CE1	2.51	0.46
2:B:33:ALA:O	2:B:37:ILE:HG13	2.16	0.46
2:B:385:LYS:HG2	2:B:386:THR:H	1.81	0.46
2:B:104:LYS:HG3	2:B:192:ASP:OD2	2.16	0.46
1:A:246:LEU:O	1:A:307:ARG:NH1	2.44	0.46
1:A:132:ILE:HB	1:A:142:ILE:CG1	2.44	0.46
1:A:218:ASP:O	1:A:222:GLN:HG3	2.16	0.46
2:B:125:ARG:NH1	2:B:147:ASN:HD22	2.13	0.46
1:A:51:GLY:C	1:A:53:GLU:H	2.18	0.46
1:A:395:LYS:CD	1:A:395:LYS:N	2.77	0.46
2:B:31:ILE:O	2:B:35:VAL:HG23	2.16	0.45
1:A:440:PHE:CE1	1:A:489:SER:HB3	2.51	0.45
2:B:270:ILE:HG12	2:B:346:PHE:HB3	1.98	0.45
2:B:5:ILE:CG2	2:B:6:GLU:N	2.78	0.45
1:A:156:SER:CB	1:A:157:PRO:HD3	2.44	0.45
2:B:401:TRP:O	2:B:404:GLU:HB2	2.17	0.45
2:B:103:LYS:HD2	2:B:191:SER:CA	2.47	0.45
2:B:168:LEU:C	2:B:170:PRO:HD2	2.36	0.45
2:B:276:VAL:O	2:B:277:ARG:C	2.55	0.45
1:A:501:TYR:CZ	1:A:505:ILE:HD11	2.51	0.45
1:A:467:VAL:CG2	1:A:484:LEU:HD11	2.47	0.44
1:A:101:LYS:H	1:A:101:LYS:CD	2.28	0.44
2:B:125:ARG:HH11	2:B:147:ASN:HD22	1.66	0.44
2:B:108:VAL:HG22	2:B:188:TYR:CD2	2.52	0.44
2:B:46:LYS:O	2:B:147:ASN:HB2	2.17	0.44
1:A:380:ILE:HG12	2:B:27:THR:HG22	1.98	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:118:VAL:HG13	1:A:119:PRO:HD2	2.00	0.44
2:B:316:GLY:O	2:B:318:TYR:HD1	2.01	0.44
2:B:111:VAL:HG22	2:B:185:ASP:O	2.18	0.44
1:A:268:SER:CB	1:A:353:LYS:HE2	2.47	0.44
1:A:366:LYS:O	1:A:370:GLU:HG3	2.18	0.44
2:B:344:GLU:O	2:B:347:LYS:HB2	2.18	0.44
1:A:232:TYR:HB3	1:A:240:THR:O	2.17	0.44
2:B:167:ILE:O	2:B:167:ILE:HG12	2.18	0.43
1:A:170:PRO:O	1:A:171:PHE:C	2.56	0.43
1:A:90:VAL:O	1:A:91:GLN:C	2.56	0.43
2:B:172:ARG:O	2:B:176:PRO:HG3	2.18	0.43
1:A:94:ILE:HG23	1:A:229:TRP:HH2	1.82	0.43
1:A:102:LYS:O	1:A:103:LYS:HD3	2.18	0.43
2:B:75:VAL:HG11	2:B:77:PHE:CZ	2.53	0.43
2:B:146:TYR:CG	2:B:150:PRO:HB3	2.53	0.43
2:B:353:LYS:HZ3	2:B:428:GLN:HG3	1.78	0.43
1:A:54:ASN:O	1:A:143:ARG:NH2	2.51	0.43
1:A:325:LEU:HD21	1:A:385:LYS:HE3	2.01	0.43
1:A:283:LEU:O	1:A:286:THR:HG23	2.18	0.43
2:B:66:LYS:HA	2:B:407:GLN:HE22	1.84	0.43
1:A:169:GLU:N	1:A:170:PRO:CD	2.81	0.43
1:A:108:VAL:C	1:A:109:LEU:HD12	2.39	0.43
2:B:160:PHE:O	2:B:160:PHE:CD1	2.72	0.43
1:A:207:GLN:O	1:A:211:ARG:HG3	2.19	0.43
1:A:153:TRP:CG	1:A:154:LYS:N	2.86	0.43
2:B:107:THR:HG22	2:B:109:LEU:CD1	2.49	0.43
1:A:41:MET:HB3	1:A:46:LYS:HG3	2.00	0.43
2:B:393:ILE:CG1	2:B:394:GLN:N	2.80	0.43
2:B:330:GLN:HB2	2:B:338:THR:HG1	1.82	0.43
1:A:88:TRP:CE3	1:A:88:TRP:HA	2.54	0.43
1:A:518:VAL:O	1:A:519:ASN:C	2.58	0.43
2:B:265:ASN:O	2:B:268:SER:OG	2.28	0.43
2:B:207:GLN:HA	2:B:210:LEU:HB2	2.01	0.42
2:B:98:ALA:O	2:B:101:LYS:HG2	2.18	0.42
1:A:93:GLY:C	1:A:94:ILE:HD12	2.39	0.42
2:B:247:PRO:O	2:B:252:TRP:CH2	2.72	0.42
1:A:246:LEU:HD12	1:A:307:ARG:CA	2.50	0.42
1:A:83:ARG:HH11	1:A:83:ARG:CG	2.27	0.42
1:A:163:SER:O	1:A:164:MET:C	2.57	0.42
1:A:320:ASP:HA	1:A:321:PRO:HD2	1.83	0.42
2:B:85:GLN:CG	2:B:87:PHE:CZ	3.02	0.42
1:A:401:TRP:O	1:A:404:GLU:HG2	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:410:TRP:CE3	2:B:363:ASN:HB2	2.55	0.42
1:A:281:LYS:O	1:A:284:ARG:HG3	2.19	0.42
2:B:8:VAL:O	2:B:121:ASP:HB2	2.19	0.42
1:A:33:ALA:HB2	1:A:71:TRP:CD1	2.54	0.42
2:B:103:LYS:NZ	2:B:177:ASP:O	2.53	0.42
1:A:328:GLU:HG3	1:A:390:LYS:CB	2.45	0.42
2:B:115:TYR:OH	2:B:157:PRO:HG3	2.20	0.42
2:B:64:LYS:HD2	2:B:70:LYS:O	2.19	0.42
1:A:442:VAL:HB	1:A:481:ALA:HB1	2.02	0.42
2:B:120:LEU:HD23	2:B:125:ARG:HG2	2.00	0.42
1:A:364:ASP:O	1:A:368:LEU:HB2	2.20	0.42
2:B:420:PRO:HA	2:B:421:PRO:HD3	1.93	0.42
1:A:60:VAL:CG2	1:A:130:PHE:HB2	2.49	0.42
1:A:278:GLN:HA	1:A:278:GLN:NE2	2.35	0.42
1:A:201:LYS:O	1:A:204:GLU:HB2	2.19	0.42
1:A:340:GLN:HE21	1:A:340:GLN:HB3	1.67	0.42
2:B:394:GLN:O	2:B:395:LYS:C	2.58	0.41
1:A:108:VAL:HG23	1:A:227:PHE:CE1	2.54	0.41
1:A:303:LEU:C	1:A:303:LEU:HD23	2.40	0.41
1:A:54:ASN:ND2	1:A:129:ALA:HB2	2.35	0.41
2:B:263:LYS:O	2:B:266:TRP:HB3	2.21	0.41
1:A:83:ARG:NH1	1:A:83:ARG:CG	2.82	0.41
1:A:479:LEU:HB3	1:A:517:LEU:HD13	2.03	0.41
2:B:393:ILE:CG1	2:B:394:GLN:H	2.32	0.41
2:B:210:LEU:HD13	2:B:210:LEU:HA	1.87	0.41
2:B:193:LEU:HB3	2:B:197:GLN:HB2	2.01	0.41
1:A:31:ILE:O	1:A:35:VAL:HG23	2.20	0.41
1:A:73:LYS:NZ	1:A:146:TYR:OH	2.54	0.41
1:A:4:PRO:HG2	1:A:212:TRP:HE3	1.85	0.41
1:A:399:GLU:HG3	1:A:402:TRP:CE3	2.56	0.41
1:A:325:LEU:HG	1:A:387:PRO:HB3	2.03	0.41
1:A:10:VAL:HG12	1:A:11:LYS:H	1.85	0.41
1:A:178:ILE:HD13	1:A:178:ILE:N	2.35	0.41
1:A:224:GLU:HA	1:A:225:PRO:HD3	1.96	0.41
1:A:65:LYS:N	1:A:70:LYS:O	2.46	0.41
1:A:295:LEU:CD2	1:A:295:LEU:N	2.83	0.41
2:B:26:LEU:HB2	2:B:31:ILE:HD11	2.03	0.41
2:B:5:ILE:CG2	2:B:6:GLU:H	2.31	0.41
2:B:234:LEU:HD23	2:B:239:TRP:HZ2	1.85	0.41
1:A:40:GLU:HG3	1:A:44:GLU:OE1	2.20	0.41
1:A:394:GLN:O	1:A:397:THR:N	2.53	0.41
2:B:153:TRP:CH2	2:B:155:GLY:HA3	2.56	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:151:GLN:HB3	2:B:185:ASP:OD1	2.21	0.41
1:A:487:GLN:HA	1:A:524:GLN:NE2	2.36	0.41
2:B:119:PRO:HA	2:B:148:VAL:HA	2.01	0.41
1:A:167:ILE:C	1:A:170:PRO:HD2	2.39	0.40
2:B:134:SER:HB2	2:B:139:THR:HB	2.02	0.40
2:B:327:ALA:O	2:B:389:PHE:HA	2.21	0.40
1:A:239:TRP:O	1:A:315:HIS:CA	2.69	0.40
1:A:5:ILE:HD13	1:A:163:SER:HB3	2.02	0.40
1:A:489:SER:HB2	1:A:493:VAL:HG13	2.03	0.40
1:A:279:LEU:HA	1:A:282:LEU:HD23	2.03	0.40
2:B:380:ILE:O	2:B:384:GLY:CA	2.69	0.40
2:B:335:GLY:O	2:B:355:ALA:HA	2.22	0.40
1:A:441:TYR:O	1:A:457:TYR:HA	2.22	0.40
1:A:292:VAL:C	1:A:293:ILE:HD12	2.41	0.40
1:A:214:LEU:HA	1:A:214:LEU:HD23	1.87	0.40
2:B:84:THR:HB	2:B:154:LYS:HE2	2.02	0.40
2:B:195:ILE:HD11	2:B:233:GLU:CG	2.52	0.40
2:B:107:THR:HG22	2:B:109:LEU:HD12	2.03	0.40
1:A:120:LEU:O	1:A:121:ASP:C	2.60	0.40
1:A:168:LEU:C	1:A:170:PRO:HD2	2.42	0.40
1:A:171:PHE:CZ	1:A:205:LEU:HB2	2.56	0.40
1:A:60:VAL:HG21	1:A:130:PHE:CD1	2.56	0.40
2:B:323:LYS:HE3	2:B:323:LYS:HB3	1.94	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	520/560 (93%)	464 (89%)	45 (9%)	11 (2%)	11	47
2	B	393/440 (89%)	346 (88%)	43 (11%)	4 (1%)	22	70
All	All	913/1000 (91%)	810 (89%)	88 (10%)	15 (2%)	14	56

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	ALA
1	A	195	ILE
1	A	402	TRP
1	A	170	PRO
1	A	230	MET
1	A	403	THR
1	A	538	ALA
2	B	232	TYR
1	A	137	ASN
1	A	243	PRO
1	A	412	PRO
2	B	193	LEU
2	B	277	ARG
2	B	395	LYS
1	A	156	SER

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	473/499 (95%)	419 (89%)	54 (11%)	8	33
2	B	366/400 (92%)	341 (93%)	25 (7%)	22	63
All	All	839/899 (93%)	760 (91%)	79 (9%)	13	44

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR
1	A	20	LYS
1	A	46	LYS
1	A	61	PHE
1	A	67	ASP
1	A	89	GLU
1	A	123	ASP
1	A	134	SER
1	A	136	ASN
1	A	139	THR
1	A	142	ILE

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Mol	Chain	Res	Type
1	A	163	SER
1	A	165	THR
1	A	168	LEU
1	A	177	ASP
1	A	185	ASP
1	A	187	LEU
1	A	194	GLU
1	A	195	ILE
1	A	206	ARG
1	A	232	TYR
1	A	243	PRO
1	A	253	THR
1	A	264	LEU
1	A	265	ASN
1	A	279	LEU
1	A	282	LEU
1	A	283	LEU
1	A	287	LYS
1	A	290	THR
1	A	295	LEU
1	A	317	VAL
1	A	325	LEU
1	A	336	GLN
1	A	340	GLN
1	A	356	ARG
1	A	362	THR
1	A	368	LEU
1	A	394	GLN
1	A	395	LYS
1	A	396	GLU
1	A	402	TRP
1	A	423	VAL
1	A	424	LYS
1	A	470	THR
1	A	471	ASP
1	A	473	THR
1	A	476	LYS
1	A	484	LEU
1	A	493	VAL
1	A	500	GLN
1	A	503	LEU
1	A	514	GLU

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Mol	Chain	Res	Type
1	A	517	LEU
2	B	8	VAL
2	B	16	MET
2	B	22	LYS
2	B	53	GLU
2	B	55	PRO
2	B	60	VAL
2	B	72	ARG
2	B	166	LYS
2	B	167	ILE
2	B	171	PHE
2	B	174	GLN
2	B	205	LEU
2	B	212	TRP
2	B	228	LEU
2	B	230	MET
2	B	280	CYS
2	B	283	LEU
2	B	286	THR
2	B	295	LEU
2	B	297	GLU
2	B	301	LEU
2	B	303	LEU
2	B	353	LYS
2	B	368	LEU
2	B	414	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	GLN
1	A	207	GLN
1	A	221	HIS
1	A	222	GLN
1	A	235	HIS
1	A	242	GLN
1	A	255	ASN
1	A	278	GLN
1	A	336	GLN
1	A	407	GLN
1	A	475	GLN
1	A	480	GLN

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Mol	Chain	Res	Type
1	A	507	GLN
1	A	509	GLN
1	A	520	GLN
2	B	147	ASN
2	B	151	GLN
2	B	161	GLN
2	B	174	GLN
2	B	175	ASN
2	B	207	GLN
2	B	242	GLN
2	B	269	GLN
2	B	278	GLN
2	B	336	GLN
2	B	348	ASN
2	B	407	GLN
2	B	428	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	CSD	A	280	1	7,7,8	7.24	3 (42%)	6,8,10	4.94	4 (66%)

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
1	CSD	A	280	1	-	1/3/6/8	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	280	CSD	O-C	18.77	1.24	1.11
1	A	280	CSD	OD1-SG	2.86	1.51	1.47
1	A	280	CSD	CA-C	2.16	1.52	1.48

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	CSD	OD2-SG-OD1	8.65	123.75	109.39
1	A	280	CSD	OD1-SG-CB	4.98	118.46	105.25
1	A	280	CSD	CA-CB-SG	4.70	117.52	110.82
1	A	280	CSD	C-CA-N	-4.59	109.25	113.83

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	280	CSD	CA-CB-SG-OD1

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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
3	NVP	A	999	-	18,23,23	1.24	3 (16%)	19,34,34	1.51	2 (10%)

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
3	NVP	A	999	-	-	0/0/6/6	0/0/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	999	NVP	C4-N3	2.24	1.37	1.32
3	A	999	NVP	C12-C11	2.22	1.41	1.36
3	A	999	NVP	C13-N14	2.15	1.36	1.32

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	999	NVP	C5-C6-C7	3.75	119.25	116.82
3	A	999	NVP	C6-C7-C2	-2.85	118.77	122.58

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	525/560 (93%)	-0.24	1 (0%) 93 54	36, 87, 141, 150	0
2	B	401/440 (91%)	-0.23	0 100 100	40, 88, 133, 149	0
All	All	926/1000 (92%)	-0.24	1 (0%) 93 63	36, 88, 136, 150	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	67	ASP	2.3

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	CSD	A	280	8/9	0.15	-0.68	78,81,85,89	0

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(\AA^2)	Q<0.9
3	NVP	A	999	20/20	0.23	1.18	58,63,77,78	0

6.5 Other polymers ⓘ

There are no such residues in this entry.