



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 21, 2018 – 01:10 AM EST

PDB ID : 1TL3
Title : Crystal structure of hiv-1 reverse transcriptase in complex with gw450557
Authors : Hopkins, A.L.; Ren, J.; Stuart, D.I.; Stammers, D.K.
Deposited on : 2004-06-09
Resolution : 2.80 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030736

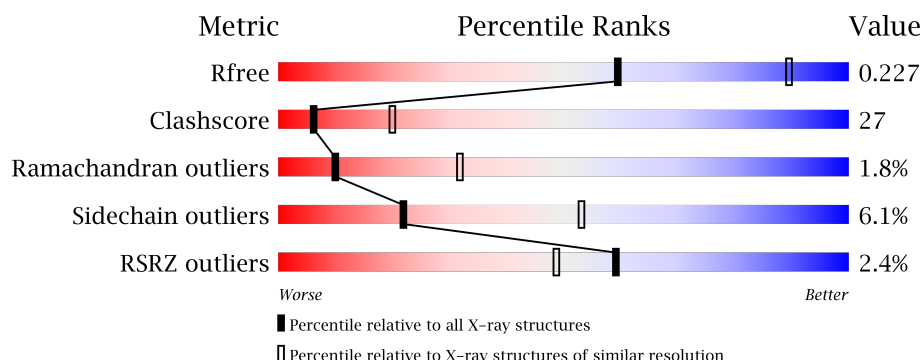
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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	560	<div> <div>2%</div> <div>46%</div> <div>41%</div> <div>5%</div> <div>8%</div> </div>
2	B	440	<div> <div>2%</div> <div>43%</div> <div>45%</div> <div>•</div> <div>9%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7612 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 Pol polyprotein, Reverse transcriptase, Chain A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	517	Total	C	N	O	S	0	0	0
			4247	2754	701	784	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	280	CSD	CYS	OXIDIZED CYS	UNP P04585

- Molecule 2 is a protein called Pol polyprotein, Reverse transcriptase, Chain B.

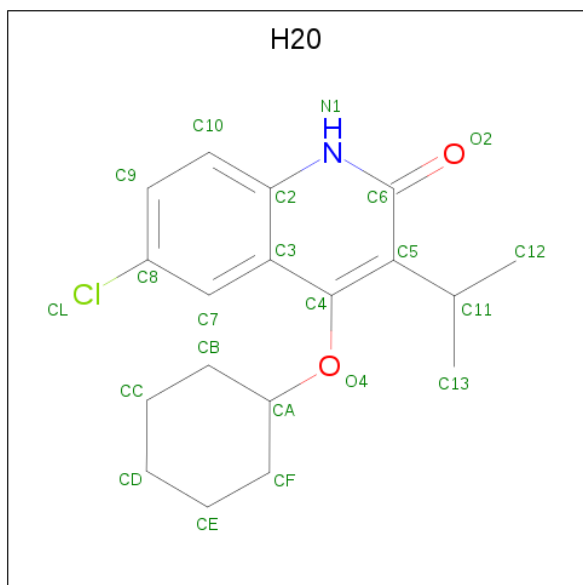
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	401	Total	C	N	O	S	0	0	0
			3328	2173	548	600	7			

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is 6-CHLORO-4-(CYCLOHEXYLOXY)-3-ISOPROPYLQUINOLIN-2(1H)-ONE (three-letter code: H20) (formula: C₁₈H₂₂ClNO₂).

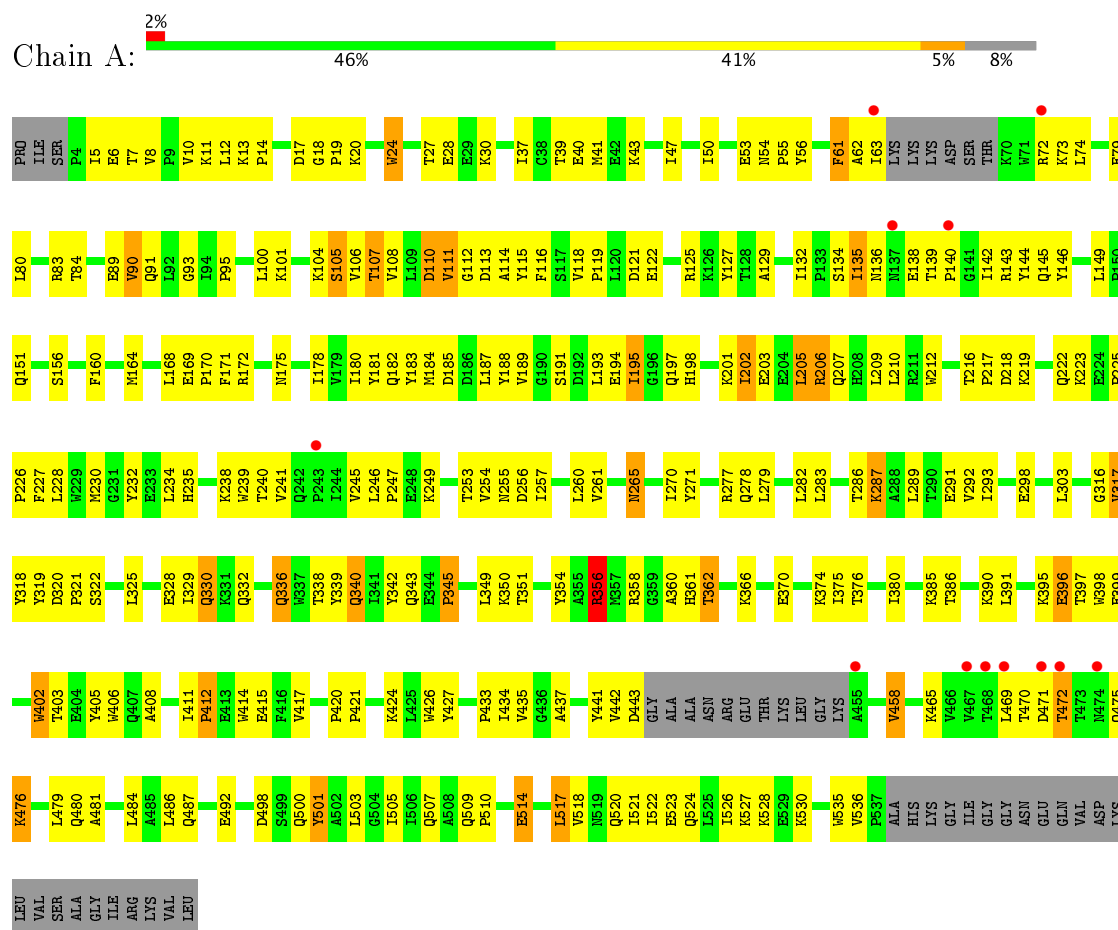


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	0	0
			22	18	1	1	2		

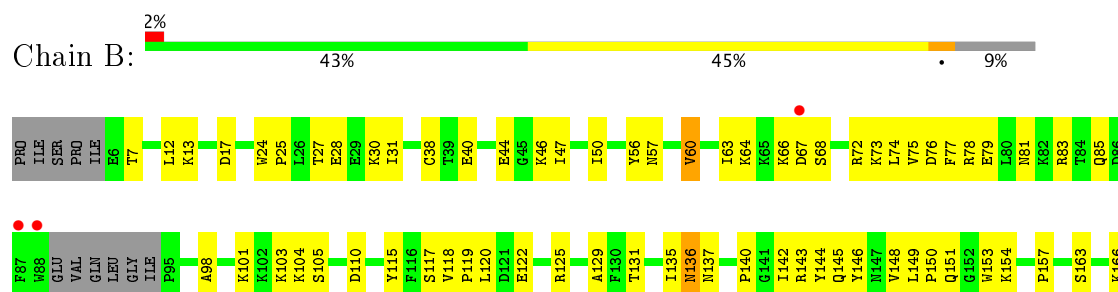
3 Residue-property plots [i](#)

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

- Molecule 1: Pol polyprotein, Reverse transcriptase, Chain A



- Molecule 2: Pol polyprotein, Reverse transcriptase, Chain B



WORLDWIDE
PDB
PROTEIN DATA BANK

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	138.00 Å 115.20 Å 65.30 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.85 – 2.80 29.84 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.5 (29.85-2.80) 92.5 (29.84-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.80 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.206 , 0.286 0.197 , 0.227	Depositor DCC
R_{free} test set	1171 reflections (4.81%)	DCC
Wilson B-factor (Å ²)	55.8	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 78.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7612	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: CSD, PO4, H2O

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	0.50	0/4352	0.71	0/5918
2	B	0.49	0/3425	0.70	0/4652
All	All	0.50	0/7777	0.70	0/10570

There are no bond length outliers.

There are no bond angle outliers.

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	4247	0	4275	234	0
2	B	3328	0	3353	202	0
3	A	15	0	0	1	0
4	A	22	0	22	2	0
All	All	7612	0	7650	414	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 27.

All (414) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:GLN:HE21	2:B:422:LEU:HD12	1.31	0.94
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.49	0.92
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.51	0.92
2:B:234:LEU:HD21	2:B:377:THR:HG21	1.52	0.91
1:A:253:THR:HG22	1:A:255:ASN:H	1.37	0.88
1:A:500:GLN:NE2	2:B:422:LEU:HD12	1.89	0.88
2:B:227:PHE:HB3	2:B:231:GLY:HA2	1.55	0.88
2:B:66:LYS:HE3	2:B:230:MET:HG3	1.58	0.86
2:B:266:TRP:O	2:B:269:GLN:HG2	1.77	0.85
2:B:57:ASN:HD22	2:B:143:ARG:NH1	1.76	0.84
1:A:354:TYR:HD2	1:A:374:LYS:HD2	1.43	0.83
2:B:362:THR:HG22	2:B:366:LYS:HD3	1.60	0.82
1:A:101:LYS:N	1:A:101:LYS:HD2	1.96	0.80
1:A:406:TRP:CH2	2:B:418:ASN:HA	2.17	0.80
2:B:193:LEU:H	2:B:193:LEU:HD12	1.49	0.77
1:A:500:GLN:HE21	2:B:422:LEU:CD1	1.97	0.77
2:B:348:ASN:HD22	2:B:351:THR:CG2	1.99	0.76
2:B:227:PHE:HB3	2:B:231:GLY:CA	2.15	0.76
1:A:469:LEU:HD21	1:A:480:GLN:HG3	1.68	0.75
2:B:203:GLU:HG3	2:B:207:GLN:NE2	2.02	0.74
1:A:91:GLN:HE21	2:B:137:ASN:HB3	1.52	0.74
1:A:206:ARG:HG2	1:A:216:THR:OG1	1.87	0.73
2:B:104:LYS:HB2	2:B:192:ASP:HA	1.70	0.73
2:B:295:LEU:HD12	2:B:295:LEU:N	2.02	0.73
2:B:193:LEU:N	2:B:193:LEU:HD12	2.03	0.73
2:B:40:GLU:HG3	2:B:44:GLU:OE2	1.89	0.73
1:A:113:ASP:HB2	1:A:116:PHE:CD2	2.24	0.73
1:A:28:GLU:OE1	1:A:135:ILE:HG22	1.89	0.71
1:A:111:VAL:HG12	1:A:114:ALA:HB2	1.71	0.71
2:B:239:TRP:HB3	2:B:350:LYS:NZ	2.07	0.70
1:A:12:LEU:HD11	1:A:127:TYR:CZ	2.27	0.69
1:A:178:ILE:HG12	1:A:191:SER:HB3	1.75	0.69
1:A:480:GLN:HG2	1:A:517:LEU:HD11	1.75	0.69
1:A:227:PHE:HB2	1:A:234:LEU:HB2	1.75	0.69
1:A:507:GLN:HE22	2:B:421:PRO:HB3	1.58	0.68
2:B:242:GLN:HG2	2:B:353:LYS:HG2	1.75	0.68
1:A:435:VAL:HG22	2:B:290:THR:HG21	1.74	0.68
2:B:366:LYS:HG3	2:B:405:TYR:CD2	2.28	0.68
1:A:412:PRO:HG3	2:B:401:TRP:CZ2	2.29	0.68
1:A:122:GLU:HA	1:A:125:ARG:HD2	1.76	0.67
1:A:239:TRP:NE1	1:A:316:GLY:HA3	2.10	0.67
1:A:354:TYR:CD2	1:A:374:LYS:HD2	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:193:LEU:H	2:B:193:LEU:CD1	2.08	0.66
1:A:249:LYS:NZ	1:A:256:ASP:HB3	2.10	0.66
1:A:206:ARG:HH21	1:A:217:PRO:C	2.00	0.66
2:B:203:GLU:HG3	2:B:207:GLN:HE22	1.61	0.66
2:B:295:LEU:HD12	2:B:295:LEU:H	1.58	0.66
2:B:110:ASP:HB3	2:B:226:PRO:HG2	1.78	0.65
1:A:111:VAL:CG1	1:A:114:ALA:HB2	2.27	0.65
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.78	0.65
2:B:254:VAL:HG23	2:B:291:GLU:O	1.97	0.65
1:A:24:TRP:HZ3	1:A:61:PHE:HB3	1.62	0.64
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.32	0.64
2:B:239:TRP:CH2	2:B:378:GLU:HG2	2.32	0.64
1:A:399:GLU:HA	1:A:402:TRP:HE3	1.61	0.64
1:A:360:ALA:HB1	1:A:514:GLU:OE1	1.97	0.64
1:A:62:ALA:C	1:A:63:ILE:HD12	2.18	0.64
1:A:91:GLN:NE2	2:B:137:ASN:HB3	2.12	0.64
1:A:228:LEU:H	1:A:228:LEU:HD12	1.61	0.64
2:B:66:LYS:HG2	2:B:230:MET:HA	1.80	0.64
1:A:206:ARG:HH21	1:A:218:ASP:HA	1.63	0.63
1:A:476:LYS:HD2	1:A:476:LYS:O	1.98	0.63
2:B:234:LEU:HD21	2:B:377:THR:CG2	2.26	0.63
1:A:507:GLN:NE2	2:B:421:PRO:HB3	2.13	0.63
1:A:476:LYS:HD3	1:A:517:LEU:HD12	1.79	0.63
2:B:183:TYR:CE1	2:B:184:MET:HG2	2.32	0.63
1:A:114:ALA:HB1	1:A:160:PHE:CE2	2.33	0.63
1:A:319:TYR:OH	1:A:385:LYS:HE2	1.98	0.63
2:B:191:SER:HB2	2:B:193:LEU:HD13	1.80	0.62
2:B:254:VAL:O	2:B:258:GLN:HG3	1.99	0.62
2:B:261:VAL:HG13	2:B:276:VAL:HG11	1.81	0.62
1:A:12:LEU:HD11	1:A:127:TYR:CE1	2.35	0.62
1:A:330:GLN:HG2	1:A:338:THR:OG1	1.99	0.62
1:A:30:LYS:HD3	1:A:62:ALA:HB3	1.81	0.62
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.82	0.62
1:A:171:PHE:CE1	1:A:205:LEU:HA	2.34	0.61
1:A:206:ARG:NH2	1:A:218:ASP:HA	2.15	0.61
1:A:105:SER:HB2	1:A:198:HIS:CD2	2.35	0.61
2:B:249:LYS:HG3	2:B:252:TRP:CE2	2.36	0.61
2:B:63:ILE:HD11	2:B:72:ARG:HD3	1.83	0.61
2:B:169:GLU:O	2:B:173:LYS:HG3	2.01	0.61
2:B:98:ALA:O	2:B:101:LYS:HG2	2.00	0.61
2:B:66:LYS:NZ	2:B:67:ASP:HB2	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.35	0.60
2:B:40:GLU:OE2	2:B:40:GLU:HA	2.01	0.60
1:A:136:ASN:OD1	1:A:139:THR:HG23	1.99	0.60
1:A:115:TYR:CD1	1:A:156:SER:HB3	2.37	0.60
1:A:95:PRO:HB3	2:B:136:ASN:O	2.01	0.60
2:B:66:LYS:HE3	2:B:230:MET:CG	2.31	0.60
1:A:249:LYS:HZ1	1:A:256:ASP:HB3	1.65	0.60
1:A:90:VAL:HG12	1:A:91:GLN:N	2.16	0.60
2:B:236:PRO:HA	2:B:239:TRP:CE2	2.36	0.60
1:A:40:GLU:HA	1:A:40:GLU:OE2	2.02	0.60
2:B:168:LEU:C	2:B:170:PRO:HD2	2.22	0.59
1:A:396:GLU:HG3	1:A:397:THR:N	2.18	0.59
1:A:342:TYR:HA	1:A:349:LEU:HD12	1.84	0.59
2:B:66:LYS:CG	2:B:230:MET:HA	2.31	0.59
1:A:125:ARG:HG2	1:A:146:TYR:O	2.03	0.59
1:A:240:THR:OG1	1:A:241:VAL:N	2.35	0.59
1:A:253:THR:HG22	1:A:255:ASN:N	2.14	0.59
1:A:24:TRP:CZ3	1:A:61:PHE:HB3	2.37	0.59
1:A:358:ARG:HD3	1:A:370:GLU:CD	2.22	0.59
1:A:206:ARG:HB3	1:A:206:ARG:HH11	1.68	0.58
1:A:522:ILE:O	1:A:526:ILE:HG13	2.04	0.58
1:A:56:TYR:O	1:A:143:ARG:NH2	2.30	0.58
2:B:7:THR:HG22	2:B:119:PRO:HG2	1.85	0.58
2:B:297:GLU:O	2:B:301:LEU:HG	2.03	0.58
1:A:277:ARG:HH11	1:A:277:ARG:HG3	1.68	0.57
2:B:193:LEU:HD22	2:B:201:LYS:HD3	1.86	0.57
2:B:282:LEU:HD21	2:B:296:THR:HG23	1.85	0.57
1:A:61:PHE:N	1:A:61:PHE:HD2	2.01	0.57
1:A:411:ILE:O	1:A:412:PRO:O	2.22	0.57
1:A:228:LEU:HA	1:A:232:TYR:O	2.05	0.57
2:B:146:TYR:CG	2:B:150:PRO:HB3	2.40	0.57
1:A:132:ILE:HB	1:A:142:ILE:HG13	1.87	0.57
2:B:174:GLN:C	2:B:176:PRO:HD3	2.25	0.56
1:A:13:LYS:HE3	1:A:84:THR:O	2.06	0.56
1:A:376:THR:HG23	1:A:386:THR:HG22	1.86	0.56
2:B:57:ASN:HD22	2:B:143:ARG:HH12	1.51	0.56
2:B:325:LEU:HD21	2:B:383:TRP:CE3	2.41	0.56
2:B:244:ILE:HD13	2:B:266:TRP:CZ3	2.41	0.56
1:A:61:PHE:N	1:A:61:PHE:CD2	2.72	0.56
1:A:18:GLY:HA3	1:A:56:TYR:CE1	2.40	0.56
1:A:279:LEU:HD23	1:A:282:LEU:HD11	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:167:ILE:O	2:B:208:HIS:HE1	1.89	0.55
2:B:166:LYS:O	2:B:169:GLU:HG3	2.06	0.55
2:B:131:THR:OG1	2:B:143:ARG:HD2	2.07	0.55
2:B:332:GLN:HA	2:B:424:LYS:HE3	1.89	0.55
1:A:246:LEU:HD22	1:A:260:LEU:HD11	1.88	0.55
1:A:115:TYR:O	1:A:149:LEU:HB2	2.06	0.55
1:A:24:TRP:HZ3	1:A:61:PHE:CG	2.25	0.55
1:A:412:PRO:HG3	2:B:401:TRP:HZ2	1.72	0.55
1:A:40:GLU:OE2	1:A:43:LYS:HD3	2.07	0.55
2:B:183:TYR:OH	2:B:386:THR:HG23	2.07	0.55
1:A:257:ILE:O	1:A:261:VAL:HG23	2.06	0.54
2:B:120:LEU:HD23	2:B:125:ARG:HG2	1.88	0.54
2:B:195:ILE:HG23	2:B:196:GLY:N	2.22	0.54
1:A:498:ASP:HA	1:A:536:VAL:O	2.08	0.54
2:B:276:VAL:O	2:B:277:ARG:C	2.45	0.54
1:A:332:GLN:HG2	1:A:332:GLN:O	2.08	0.54
2:B:298:GLU:CD	2:B:298:GLU:H	2.11	0.54
2:B:197:GLN:O	2:B:201:LYS:HG3	2.07	0.54
1:A:228:LEU:N	1:A:228:LEU:HD12	2.22	0.54
1:A:216:THR:HB	1:A:217:PRO:HD2	1.90	0.53
2:B:125:ARG:HB3	2:B:145:GLN:NE2	2.23	0.53
1:A:265:ASN:N	1:A:265:ASN:HD22	2.05	0.53
1:A:329:ILE:HD11	1:A:375:ILE:HD12	1.90	0.53
2:B:305:GLU:O	2:B:309:ILE:HG13	2.09	0.53
2:B:295:LEU:CD1	2:B:295:LEU:N	2.72	0.53
1:A:406:TRP:O	2:B:331:LYS:HB3	2.08	0.53
2:B:365:VAL:HG11	2:B:401:TRP:HB2	1.90	0.53
2:B:182:GLN:HA	2:B:187:LEU:HD12	1.91	0.53
1:A:469:LEU:HD21	1:A:480:GLN:CG	2.38	0.53
2:B:66:LYS:HZ2	2:B:67:ASP:HB2	1.72	0.53
1:A:441:TYR:HB2	1:A:458:VAL:HG23	1.90	0.53
2:B:199:ARG:CZ	2:B:233:GLU:OE1	2.56	0.53
1:A:406:TRP:CZ3	2:B:418:ASN:HA	2.44	0.53
1:A:116:PHE:CD2	1:A:116:PHE:N	2.77	0.53
1:A:498:ASP:O	1:A:535:TRP:NE1	2.41	0.53
1:A:442:VAL:HB	1:A:481:ALA:HB1	1.91	0.52
2:B:63:ILE:HG21	2:B:74:LEU:HD22	1.91	0.52
2:B:66:LYS:HD2	2:B:67:ASP:N	2.25	0.52
2:B:195:ILE:CG2	2:B:196:GLY:N	2.73	0.52
1:A:399:GLU:HG3	1:A:402:TRP:CE3	2.45	0.52
2:B:208:HIS:HA	2:B:211:ARG:HD3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:521:ILE:O	1:A:524:GLN:HB2	2.09	0.52
1:A:5:ILE:HG13	1:A:6:GLU:N	2.23	0.52
1:A:362:THR:HG22	1:A:366:LYS:HD3	1.93	0.51
1:A:505:ILE:O	1:A:510:PRO:HD3	2.10	0.51
2:B:229:TRP:CE3	2:B:229:TRP:HA	2.46	0.51
2:B:64:LYS:HE2	2:B:68:SER:O	2.10	0.51
2:B:202:ILE:HG21	2:B:227:PHE:HE1	1.74	0.51
1:A:37:ILE:O	1:A:41:MET:HG3	2.11	0.51
2:B:175:ASN:N	2:B:176:PRO:HD3	2.25	0.51
1:A:398:TRP:CE2	1:A:411:ILE:HD12	2.45	0.51
1:A:292:VAL:C	1:A:293:ILE:HD12	2.31	0.51
1:A:91:GLN:C	1:A:93:GLY:H	2.14	0.51
2:B:239:TRP:HB3	2:B:350:LYS:HZ1	1.74	0.51
2:B:239:TRP:CZ3	2:B:378:GLU:HG2	2.46	0.51
2:B:380:ILE:O	2:B:384:GLY:N	2.43	0.51
1:A:142:ILE:HD12	1:A:144:TYR:OH	2.11	0.51
1:A:101:LYS:HD2	1:A:101:LYS:H	1.73	0.51
1:A:175:ASN:OD1	1:A:201:LYS:HE3	2.11	0.51
1:A:129:ALA:HA	1:A:144:TYR:O	2.11	0.51
1:A:210:LEU:C	1:A:212:TRP:N	2.63	0.50
1:A:30:LYS:CD	1:A:62:ALA:HB3	2.40	0.50
2:B:195:ILE:HD11	2:B:233:GLU:HG3	1.93	0.50
2:B:330:GLN:HB2	2:B:338:THR:OG1	2.10	0.50
1:A:476:LYS:CD	1:A:517:LEU:HD12	2.41	0.50
2:B:278:GLN:NE2	2:B:298:GLU:HB2	2.26	0.50
2:B:57:ASN:HD22	2:B:143:ARG:HH11	1.58	0.50
2:B:295:LEU:H	2:B:295:LEU:CD1	2.22	0.50
2:B:203:GLU:CG	2:B:207:GLN:HE22	2.24	0.50
1:A:472:THR:CB	1:A:476:LYS:HE3	2.42	0.50
1:A:63:ILE:N	1:A:63:ILE:HD12	2.27	0.50
1:A:116:PHE:HE1	3:A:1302:PO4:O1	1.94	0.49
1:A:18:GLY:HA3	1:A:56:TYR:CD1	2.47	0.49
2:B:348:ASN:HD22	2:B:351:THR:HG21	1.76	0.49
1:A:107:THR:HG21	1:A:222:GLN:NE2	2.27	0.49
1:A:226:PRO:HA	1:A:234:LEU:O	2.12	0.49
2:B:142:ILE:HG22	2:B:144:TYR:CE2	2.48	0.49
2:B:339:TYR:CD1	2:B:375:ILE:HD11	2.47	0.49
1:A:279:LEU:HA	1:A:282:LEU:CD1	2.43	0.49
1:A:61:PHE:CE2	1:A:74:LEU:HG	2.47	0.49
2:B:182:GLN:HB2	2:B:187:LEU:CD1	2.42	0.49
1:A:350:LYS:HG2	1:A:351:THR:N	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:ILE:HD13	2:B:27:THR:HG22	1.95	0.49
1:A:225:PRO:HG3	1:A:227:PHE:CE2	2.47	0.49
1:A:492:GLU:HG2	1:A:530:LYS:HB2	1.94	0.49
2:B:125:ARG:HB3	2:B:145:GLN:HE21	1.76	0.49
2:B:12:LEU:HD22	2:B:83:ARG:O	2.13	0.49
2:B:342:TYR:HB3	2:B:348:ASN:HA	1.94	0.49
1:A:391:LEU:C	1:A:417:VAL:HG12	2.33	0.49
1:A:472:THR:OG1	1:A:476:LYS:HE3	2.13	0.49
2:B:151:GLN:HB3	2:B:185:ASP:OD1	2.11	0.49
2:B:393:ILE:HG12	2:B:394:GLN:N	2.28	0.49
1:A:210:LEU:C	1:A:212:TRP:H	2.16	0.49
1:A:47:ILE:HD12	1:A:144:TYR:CD2	2.48	0.48
1:A:106:VAL:HA	1:A:189:VAL:O	2.13	0.48
2:B:146:TYR:CD2	2:B:150:PRO:HB3	2.48	0.48
1:A:426:TRP:N	1:A:426:TRP:CD1	2.81	0.48
2:B:335:GLY:O	2:B:355:ALA:HA	2.14	0.48
1:A:169:GLU:O	1:A:172:ARG:HB2	2.14	0.48
1:A:376:THR:HG23	1:A:386:THR:CG2	2.43	0.48
1:A:329:ILE:HG22	1:A:330:GLN:N	2.29	0.48
1:A:470:THR:O	1:A:471:ASP:HB2	2.13	0.48
1:A:486:LEU:O	1:A:528:LYS:NZ	2.45	0.48
1:A:523:GLU:O	1:A:527:LYS:HG2	2.14	0.48
2:B:169:GLU:N	2:B:170:PRO:HD2	2.28	0.48
2:B:191:SER:HB2	2:B:193:LEU:CD1	2.43	0.48
2:B:66:LYS:HG2	2:B:230:MET:CA	2.42	0.48
1:A:253:THR:CG2	1:A:289:LEU:O	2.62	0.48
1:A:278:GLN:HA	1:A:278:GLN:NE2	2.29	0.48
2:B:50:ILE:CG2	2:B:145:GLN:HG2	2.43	0.48
2:B:154:LYS:O	2:B:157:PRO:HD2	2.14	0.48
2:B:173:LYS:O	2:B:176:PRO:HD3	2.14	0.48
1:A:408:ALA:HB2	2:B:337:TRP:HH2	1.79	0.47
1:A:358:ARG:NH2	2:B:394:GLN:HG2	2.29	0.47
1:A:89:GLU:O	1:A:90:VAL:C	2.52	0.47
2:B:31:ILE:HD12	2:B:135:ILE:HD11	1.96	0.47
2:B:205:LEU:O	2:B:209:LEU:HG	2.14	0.47
1:A:235:HIS:HB2	1:A:238:LYS:O	2.14	0.47
1:A:492:GLU:OE1	1:A:530:LYS:HD2	2.14	0.47
2:B:257:ILE:O	2:B:260:LEU:HB3	2.14	0.47
1:A:340:GLN:HA	1:A:351:THR:HA	1.96	0.47
1:A:395:LYS:HD3	1:A:414:TRP:CZ2	2.49	0.47
2:B:270:ILE:O	2:B:272:PRO:HD3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:LEU:HD23	1:A:343:GLN:HG2	1.95	0.47
2:B:117:SER:O	2:B:118:VAL:CG2	2.63	0.47
2:B:12:LEU:O	2:B:13:LYS:C	2.51	0.47
2:B:105:SER:O	2:B:190:GLY:HA2	2.15	0.47
1:A:434:ILE:HD13	1:A:530:LYS:HB3	1.97	0.47
2:B:376:THR:O	2:B:377:THR:C	2.54	0.47
1:A:206:ARG:NH2	1:A:217:PRO:O	2.47	0.46
1:A:469:LEU:CD2	1:A:480:GLN:HG3	2.39	0.46
2:B:163:SER:O	2:B:167:ILE:HG13	2.16	0.46
2:B:397:THR:O	2:B:400:THR:HB	2.15	0.46
2:B:57:ASN:ND2	2:B:143:ARG:NH1	2.55	0.46
2:B:229:TRP:HZ3	2:B:407:GLN:HB2	1.81	0.46
2:B:252:TRP:CZ3	2:B:260:LEU:HD22	2.49	0.46
1:A:254:VAL:HG22	1:A:293:ILE:CD1	2.46	0.46
1:A:517:LEU:HA	1:A:520:GLN:NE2	2.30	0.46
1:A:194:GLU:O	1:A:195:ILE:C	2.53	0.46
1:A:402:TRP:CG	1:A:403:THR:N	2.83	0.46
2:B:244:ILE:HD13	2:B:266:TRP:HZ3	1.79	0.46
2:B:27:THR:OG1	2:B:30:LYS:HG2	2.16	0.46
2:B:328:GLU:HG2	2:B:390:LYS:HD2	1.97	0.46
1:A:202:ILE:HG22	1:A:203:GLU:N	2.31	0.46
1:A:218:ASP:O	1:A:222:GLN:HG3	2.16	0.46
4:A:999:H20:HB1	4:A:999:H20:H122	1.98	0.46
2:B:79:GLU:O	2:B:83:ARG:HG3	2.16	0.46
1:A:180:ILE:HA	1:A:188:TYR:O	2.16	0.46
1:A:253:THR:HB	1:A:256:ASP:OD2	2.16	0.46
2:B:167:ILE:HG23	2:B:212:TRP:CE3	2.51	0.45
1:A:181:TYR:CE1	1:A:183:TYR:HB2	2.50	0.45
1:A:24:TRP:HZ3	1:A:61:PHE:CB	2.26	0.45
1:A:254:VAL:HG23	1:A:291:GLU:O	2.15	0.45
2:B:129:ALA:HA	2:B:144:TYR:O	2.16	0.45
2:B:24:TRP:CD1	2:B:25:PRO:HD2	2.51	0.45
2:B:332:GLN:HB2	2:B:336:GLN:O	2.17	0.45
2:B:189:VAL:HG11	2:B:202:ILE:HD13	1.98	0.45
2:B:193:LEU:CD2	2:B:201:LYS:HD2	2.46	0.45
1:A:116:PHE:HD2	1:A:116:PHE:H	1.63	0.45
1:A:270:ILE:HG23	1:A:271:TYR:N	2.30	0.45
1:A:417:VAL:O	1:A:417:VAL:HG13	2.17	0.45
2:B:193:LEU:CD2	2:B:201:LYS:CD	2.95	0.45
1:A:79:GLU:OE2	1:A:83:ARG:NH1	2.48	0.45
1:A:424:LYS:NZ	1:A:426:TRP:CD2	2.76	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:PRO:HB3	2:B:289:LEU:HD23	1.98	0.45
1:A:283:LEU:O	1:A:286:THR:HG23	2.17	0.45
1:A:193:LEU:HD13	1:A:197:GLN:HG3	1.98	0.45
1:A:72:ARG:HG2	1:A:73:LYS:N	2.32	0.45
2:B:17:ASP:OD1	2:B:56:TYR:OH	2.27	0.45
1:A:408:ALA:HB1	2:B:364:ASP:HB3	1.97	0.45
1:A:328:GLU:O	1:A:339:TYR:HA	2.17	0.44
1:A:90:VAL:HG12	1:A:91:GLN:H	1.79	0.44
2:B:183:TYR:CD1	2:B:184:MET:HG2	2.51	0.44
2:B:193:LEU:HD22	2:B:201:LYS:CD	2.46	0.44
1:A:317:VAL:HG13	1:A:318:TYR:N	2.32	0.44
1:A:420:PRO:HA	1:A:421:PRO:C	2.37	0.44
1:A:479:LEU:HD11	1:A:501:TYR:CE2	2.52	0.44
2:B:208:HIS:NE2	2:B:212:TRP:HZ3	2.15	0.44
2:B:229:TRP:CZ3	2:B:407:GLN:HB2	2.53	0.44
1:A:13:LYS:HB3	1:A:14:PRO:HD2	1.99	0.44
1:A:398:TRP:NE1	1:A:411:ILE:HD12	2.32	0.44
1:A:501:TYR:CZ	1:A:505:ILE:HD11	2.53	0.44
1:A:520:GLN:HE21	1:A:520:GLN:HB2	1.61	0.44
2:B:81:ASN:OD1	2:B:153:TRP:HD1	2.00	0.44
2:B:66:LYS:CE	2:B:230:MET:HG3	2.37	0.44
2:B:303:LEU:O	2:B:306:ASN:HB2	2.18	0.44
1:A:164:MET:O	1:A:168:LEU:HG	2.18	0.44
1:A:110:ASP:O	1:A:112:GLY:N	2.51	0.43
1:A:287:LYS:HG3	1:A:287:LYS:H	1.72	0.43
1:A:361:HIS:O	1:A:362:THR:HG23	2.18	0.43
2:B:297:GLU:HA	2:B:300:GLU:HB2	2.00	0.43
2:B:38:CYS:O	2:B:47:ILE:HD11	2.18	0.43
2:B:103:LYS:HE3	2:B:192:ASP:OD1	2.18	0.43
2:B:332:GLN:HG3	2:B:338:THR:HG23	2.00	0.43
2:B:135:ILE:HG22	2:B:136:ASN:N	2.33	0.43
1:A:134:SER:HB3	1:A:140:PRO:O	2.18	0.43
1:A:279:LEU:HA	1:A:282:LEU:HD12	1.99	0.43
1:A:406:TRP:HH2	2:B:418:ASN:OD1	2.01	0.43
2:B:27:THR:O	2:B:28:GLU:C	2.57	0.43
2:B:257:ILE:HG21	2:B:283:LEU:HD13	2.00	0.43
1:A:197:GLN:HE21	1:A:201:LYS:HG2	1.84	0.43
1:A:245:VAL:HG23	1:A:245:VAL:O	2.18	0.43
1:A:61:PHE:HD2	1:A:61:PHE:H	1.66	0.43
2:B:169:GLU:OE2	2:B:170:PRO:HD3	2.18	0.43
1:A:356:ARG:CZ	1:A:358:ARG:HG2	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:187:LEU:HA	2:B:187:LEU:HD12	1.91	0.43
1:A:10:VAL:HG12	1:A:11:LYS:N	2.34	0.43
1:A:108:VAL:HA	1:A:187:LEU:O	2.19	0.43
2:B:76:ASP:C	2:B:78:ARG:H	2.21	0.43
1:A:187:LEU:HA	1:A:187:LEU:HD12	1.87	0.43
1:A:403:THR:CG2	2:B:334:GLN:H	2.32	0.43
2:B:170:PRO:HG2	2:B:171:PHE:H	1.82	0.43
1:A:210:LEU:O	1:A:212:TRP:N	2.51	0.43
1:A:277:ARG:NH1	1:A:277:ARG:HG3	2.32	0.42
1:A:390:LYS:HD2	1:A:415:GLU:OE2	2.18	0.42
1:A:91:GLN:OE1	2:B:140:PRO:HA	2.19	0.42
2:B:320:ASP:O	2:B:343:GLN:NE2	2.43	0.42
1:A:17:ASP:O	1:A:83:ARG:HD3	2.19	0.42
1:A:206:ARG:CB	1:A:206:ARG:HH11	2.32	0.42
1:A:340:GLN:HB3	1:A:340:GLN:HE21	1.61	0.42
1:A:61:PHE:CD2	1:A:74:LEU:O	2.72	0.42
2:B:142:ILE:CG2	2:B:144:TYR:CE2	3.02	0.42
1:A:207:GLN:HA	1:A:207:GLN:NE2	2.34	0.42
1:A:254:VAL:HG22	1:A:293:ILE:HD11	2.02	0.42
2:B:232:TYR:HA	2:B:232:TYR:HD2	1.72	0.42
2:B:282:LEU:HB3	2:B:293:ILE:HG21	2.02	0.42
2:B:306:ASN:HA	2:B:309:ILE:HD12	2.02	0.42
2:B:327:ALA:O	2:B:389:PHE:HA	2.20	0.42
2:B:425:LEU:HD23	2:B:425:LEU:O	2.19	0.42
1:A:226:PRO:HG3	1:A:235:HIS:CE1	2.54	0.42
2:B:207:GLN:HA	2:B:210:LEU:HD12	2.02	0.42
1:A:465:LYS:HD2	1:A:484:LEU:HD22	2.02	0.42
1:A:53:GLU:O	1:A:55:PRO:HD3	2.20	0.42
1:A:50:ILE:CG2	1:A:145:GLN:HG2	2.49	0.42
1:A:320:ASP:OD2	1:A:322:SER:OG	2.35	0.42
1:A:332:GLN:HB3	1:A:336:GLN:HB3	2.02	0.42
1:A:500:GLN:O	1:A:503:LEU:HB3	2.19	0.42
2:B:117:SER:C	2:B:118:VAL:HG23	2.40	0.42
1:A:108:VAL:HG13	1:A:223:LYS:HB2	2.01	0.42
1:A:188:TYR:CD2	4:A:999:H20:HF1	2.55	0.42
2:B:241:VAL:O	2:B:243:PRO:HD3	2.20	0.42
1:A:19:PRO:HD3	1:A:80:LEU:HD13	2.02	0.41
2:B:379:SER:CB	2:B:387:PRO:HD3	2.50	0.41
2:B:387:PRO:HG2	2:B:389:PHE:CE1	2.55	0.41
1:A:114:ALA:CB	1:A:160:PHE:CE2	3.02	0.41
1:A:319:TYR:O	1:A:321:PRO:HD3	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:518:VAL:O	1:A:522:ILE:HG13	2.20	0.41
2:B:117:SER:O	2:B:118:VAL:HG23	2.20	0.41
2:B:66:LYS:HG3	2:B:230:MET:HA	2.02	0.41
2:B:284:ARG:O	2:B:287:LYS:NZ	2.47	0.41
1:A:8:VAL:O	1:A:121:ASP:HB2	2.20	0.41
1:A:427:TYR:CE2	1:A:509:GLN:HG2	2.55	0.41
1:A:50:ILE:HD12	1:A:54:ASN:CB	2.51	0.41
2:B:146:TYR:CE2	2:B:150:PRO:HA	2.56	0.41
1:A:100:LEU:O	1:A:318:TYR:HB3	2.21	0.41
2:B:203:GLU:CD	2:B:207:GLN:HE22	2.24	0.41
2:B:380:ILE:O	2:B:381:VAL:C	2.59	0.41
2:B:229:TRP:HA	2:B:229:TRP:HE3	1.82	0.41
1:A:434:ILE:HB	1:A:437:ALA:HB3	2.02	0.41
2:B:372:VAL:HG13	2:B:389:PHE:CZ	2.55	0.41
2:B:325:LEU:O	2:B:387:PRO:HA	2.20	0.41
1:A:104:LYS:HB2	1:A:104:LYS:HE3	1.91	0.41
1:A:183:TYR:HE1	1:A:184:MET:CE	2.34	0.41
2:B:367:GLN:HA	2:B:370:GLU:OE1	2.21	0.41
2:B:368:LEU:O	2:B:372:VAL:HG23	2.21	0.41
1:A:239:TRP:CD1	1:A:316:GLY:C	2.94	0.41
1:A:255:ASN:HB2	1:A:289:LEU:O	2.21	0.41
1:A:405:TYR:O	2:B:331:LYS:HD3	2.21	0.41
2:B:285:GLY:O	2:B:287:LYS:N	2.54	0.41
2:B:317:VAL:O	2:B:317:VAL:HG12	2.21	0.41
1:A:27:THR:O	1:A:30:LYS:N	2.53	0.41
2:B:228:LEU:HD23	2:B:228:LEU:HA	1.79	0.41
2:B:282:LEU:CD2	2:B:296:THR:HG23	2.51	0.41
1:A:219:LYS:HD2	1:A:222:GLN:NE2	2.34	0.40
1:A:83:ARG:HG3	1:A:83:ARG:HH11	1.86	0.40
2:B:47:ILE:HD12	2:B:144:TYR:CD1	2.56	0.40
1:A:486:LEU:HD13	1:A:524:GLN:HB2	2.03	0.40
2:B:46:LYS:HA	2:B:148:VAL:HG13	2.03	0.40
2:B:328:GLU:CG	2:B:390:LYS:HD2	2.51	0.40
2:B:85:GLN:O	2:B:85:GLN:HG3	2.22	0.40
1:A:91:GLN:C	1:A:93:GLY:N	2.75	0.40
2:B:303:LEU:HD22	2:B:307:ARG:HG3	2.04	0.40
1:A:118:VAL:HG13	1:A:119:PRO:HD2	2.02	0.40
1:A:205:LEU:HD22	1:A:209:LEU:HG	2.03	0.40
1:A:246:LEU:HA	1:A:247:PRO:HD3	1.98	0.40
2:B:203:GLU:OE2	2:B:207:GLN:NE2	2.48	0.40
1:A:206:ARG:HH21	1:A:218:ASP:CA	2.31	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:GLU:OE2	1:A:298:GLU:N	2.50	0.40
1:A:486:LEU:CD1	1:A:521:ILE:HG23	2.51	0.40
2:B:195:ILE:HG13	2:B:199:ARG:HE	1.87	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	510/560 (91%)	454 (89%)	45 (9%)	11 (2%)	8	26
2	B	393/440 (89%)	340 (86%)	48 (12%)	5 (1%)	14	41
All	All	903/1000 (90%)	794 (88%)	93 (10%)	16 (2%)	10	32

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	VAL
1	A	111	VAL
1	A	195	ILE
1	A	402	TRP
1	A	412	PRO
1	A	472	THR
2	B	136	ASN
1	A	230	MET
1	A	345	PRO
2	B	77	PHE
2	B	277	ARG
1	A	356	ARG
2	B	286	THR
1	A	501	TYR
2	B	250	ASP

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Mol	Chain	Res	Type
1	A	170	PRO

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	466/499 (93%)	433 (93%)	33 (7%)	17	44
2	B	366/400 (92%)	348 (95%)	18 (5%)	29	62
All	All	832/899 (92%)	781 (94%)	51 (6%)	22	53

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

Mol	Chain	Res	Type
1	A	7	THR
1	A	20	LYS
1	A	24	TRP
1	A	39	THR
1	A	61	PHE
1	A	105	SER
1	A	107	THR
1	A	110	ASP
1	A	135	ILE
1	A	138	GLU
1	A	151	GLN
1	A	182	GLN
1	A	185	ASP
1	A	202	ILE
1	A	205	LEU
1	A	206	ARG
1	A	265	ASN
1	A	287	LYS
1	A	303	LEU
1	A	317	VAL
1	A	330	GLN
1	A	336	GLN

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Mol	Chain	Res	Type
1	A	340	GLN
1	A	345	PRO
1	A	356	ARG
1	A	362	THR
1	A	396	GLU
1	A	443	ASP
1	A	458	VAL
1	A	476	LYS
1	A	487	GLN
1	A	514	GLU
1	A	517	LEU
2	B	60	VAL
2	B	73	LYS
2	B	122	GLU
2	B	205	LEU
2	B	206	ARG
2	B	212	TRP
2	B	232	TYR
2	B	233	GLU
2	B	240	THR
2	B	249	LYS
2	B	250	ASP
2	B	295	LEU
2	B	298	GLU
2	B	303	LEU
2	B	405	TYR
2	B	410	TRP
2	B	413	GLU
2	B	414	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	197	GLN
1	A	207	GLN
1	A	208	HIS
1	A	222	GLN
1	A	265	ASN
1	A	278	GLN
1	A	330	GLN
1	A	336	GLN

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Mol	Chain	Res	Type
1	A	373	GLN
1	A	407	GLN
1	A	475	GLN
1	A	480	GLN
1	A	500	GLN
1	A	507	GLN
1	A	512	GLN
1	A	520	GLN
2	B	57	ASN
2	B	137	ASN
2	B	147	ASN
2	B	278	GLN
2	B	336	GLN
2	B	348	ASN
2	B	394	GLN
2	B	407	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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	4,7,8	1.35	1 (25%)	2,8,10	4.63	1 (50%)

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	-	1/2/6/8	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	280	CSD	CA-C	2.34	1.53	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	CSD	OD1-SG-CB	6.45	117.69	105.61

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.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

4 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	$\# Z > 2$
3	PO4	A	1300	-	4,4,4	1.29	0	6,6,6	0.39	0
3	PO4	A	1301	-	4,4,4	1.30	0	6,6,6	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PO4	A	1302	-	4,4,4	1.35	0	6,6,6	0.39	0
4	H2O	A	999	-	23,24,24	2.22	8 (34%)	28,34,34	1.95	6 (21%)

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	PO4	A	1300	-	-	0/0/0/0	0/0/0/0
3	PO4	A	1301	-	-	0/0/0/0	0/0/0/0
3	PO4	A	1302	-	-	0/0/0/0	0/0/0/0
4	H2O	A	999	-	-	0/8/16/16	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	999	H2O	C8-CL	-2.81	1.68	1.74
4	A	999	H2O	CE-CF	2.23	1.59	1.53
4	A	999	H2O	C10-C9	2.48	1.41	1.36
4	A	999	H2O	O4-C4	3.24	1.45	1.38
4	A	999	H2O	C6-N1	3.46	1.39	1.33
4	A	999	H2O	C4-C5	3.72	1.41	1.36
4	A	999	H2O	CB-CA	4.10	1.61	1.51
4	A	999	H2O	CF-CA	5.53	1.65	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	999	H2O	C3-C2-N1	-2.94	120.60	123.38
4	A	999	H2O	C6-C5-C11	-2.12	116.27	122.80
4	A	999	H2O	O4-CA-CF	2.40	113.63	108.34
4	A	999	H2O	CE-CF-CA	2.49	116.90	110.56
4	A	999	H2O	O4-CA-CB	4.23	117.67	108.34
4	A	999	H2O	C6-N1-C2	6.34	121.38	116.42

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1302	PO4	1	0
4	A	999	H2O	2	0

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	516/560 (92%)	-0.33	12 (2%)	61 51	20, 58, 100, 131	0
2	B	401/440 (91%)	-0.19	10 (2%)	58 47	18, 57, 104, 126	0
All	All	917/1000 (91%)	-0.27	22 (2%)	59 49	18, 58, 101, 131	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	469	LEU	4.2
2	B	88	TRP	4.0
1	A	468	THR	3.2
2	B	294	PRO	3.1
1	A	72	ARG	3.1
2	B	362	THR	2.9
1	A	137	ASN	2.9
1	A	243	PRO	2.7
2	B	214	LEU	2.6
2	B	237	ASP	2.6
1	A	63	ILE	2.5
1	A	467	VAL	2.5
1	A	140	PRO	2.4
2	B	87	PHE	2.3
2	B	67	ASP	2.2
2	B	356	ARG	2.2
2	B	212	TRP	2.1
2	B	334	GLN	2.1
1	A	474	ASN	2.1
1	A	455	ALA	2.1
1	A	471	ASP	2.0
1	A	472	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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
1	CSD	A	280	8/9	0.95	0.14	-	50,52,65,75	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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
3	PO4	A	1301	5/5	0.91	0.15	0.85	117,124,127,128	0
3	PO4	A	1302	5/5	0.95	0.22	0.70	134,137,138,141	0
3	PO4	A	1300	5/5	0.88	0.18	0.64	113,116,121,122	0
4	H2O	A	999	22/22	0.97	0.17	0.48	24,41,52,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.