



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:14 AM UTC

PDB ID : 2DU5 / pdb_00002du5
Title : Crystal structure of Archaeoglobus fulgidus O-phosphoseryl-tRNA synthetase E418N/E420N mutant complexed with tRNA^{Opal} and O-phosphoserine ("opal complex")
Authors : Fukunaga, R.
Deposited on : 2006-07-20
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

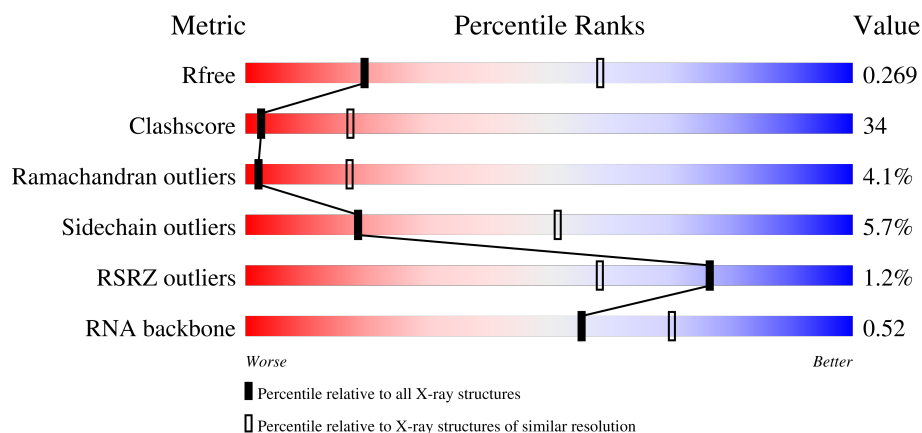
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)
RNA backbone	3983	1222 (3.50-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	D	71	 49% 34% 15% 2%
2	A	534	 41% 52% 7% 2%
2	B	534	 37% 44% 6% 13%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SEP	A	1001	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9619 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 RNA chain called tRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	71	Total	C	N	O	P	0	0	0
			1517	674	269	503	71			

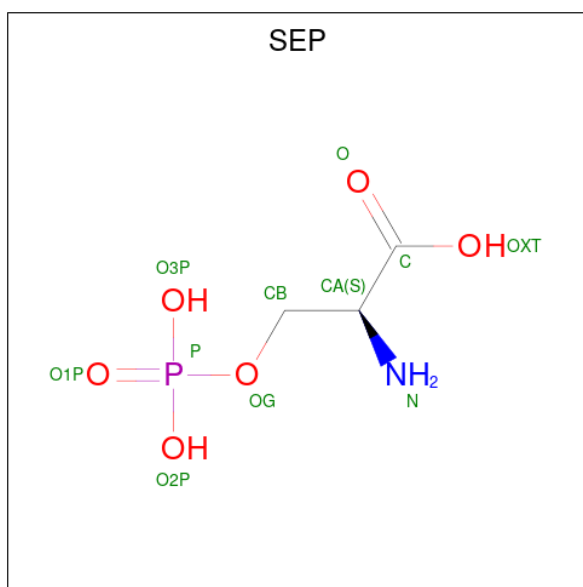
- Molecule 2 is a protein called O-phosphoseryl-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	534	Total	C	N	O	S	0	0	0
			4319	2784	721	801	13			
2	B	465	Total	C	N	O	S	0	0	0
			3761	2427	632	689	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	418	ASN	GLU	engineered mutation	UNP O30126
A	420	ASN	GLU	engineered mutation	UNP O30126
B	418	ASN	GLU	engineered mutation	UNP O30126
B	420	ASN	GLU	engineered mutation	UNP O30126

- Molecule 3 is PHOSPHOSERINE (CCD ID: SEP) (formula: C₃H₈NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	11	3	1	6	1	0	0
3	A	1	Total	C	N	O	P	0	0
			11	3	1	6	1		

3 Residue-property plots

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

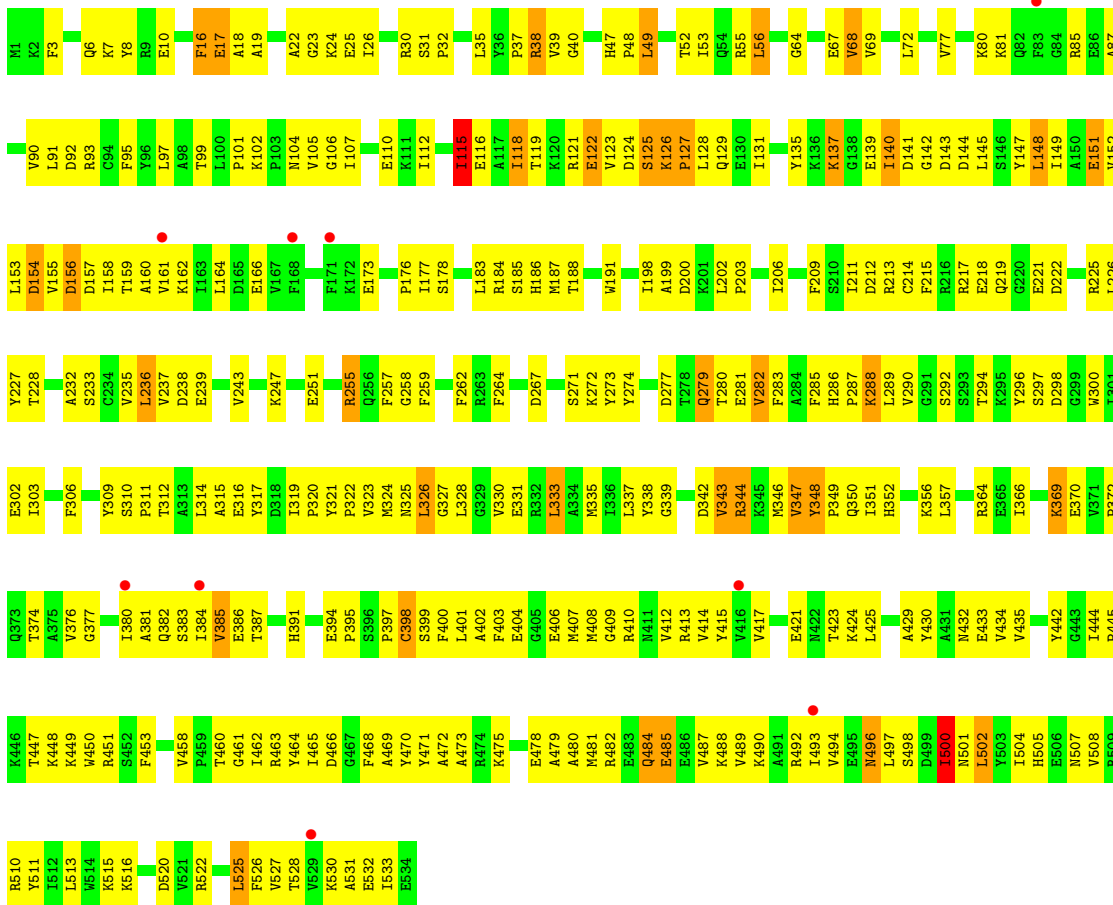
• Molecule 1: tRNA

Chain D: 

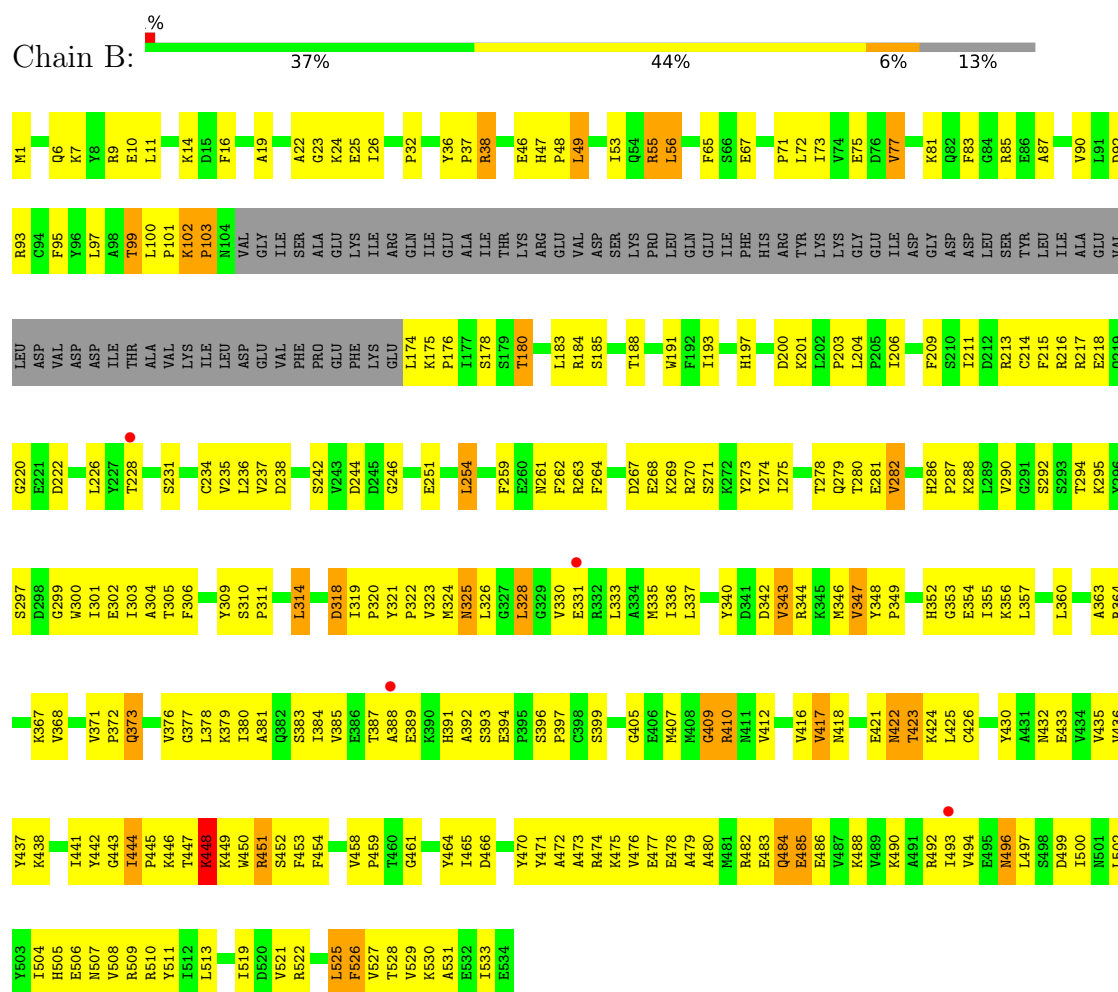


• Molecule 2: O-phosphoseryl-tRNA synthetase

Chain A: 



• Molecule 2: O-phosphoseryl-tRNA synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	148.04Å 148.04Å 152.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.07 – 3.20 49.07 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.07-3.20) 99.8 (49.07-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 3.19Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.299 0.208 , 0.269	Depositor DCC
R_{free} test set	1603 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	102.1	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 118.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9619	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% 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: SEP

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	D	0.31	0/1693	0.61	1/2637 (0.0%)
2	A	0.36	0/4416	0.86	9/5966 (0.2%)
2	B	0.39	0/3849	0.85	6/5200 (0.1%)
All	All	0.37	0/9958	0.82	16/13803 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	319	ILE	CA-C-N	8.80	128.52	119.19
2	A	319	ILE	C-N-CA	8.80	128.52	119.19
2	B	278	THR	N-CA-C	-6.10	106.38	113.88
2	B	444	ILE	N-CA-C	5.96	115.75	109.01
2	B	461	GLY	N-CA-C	-5.82	107.29	115.32
2	A	385	VAL	N-CA-C	-5.61	105.64	110.74
2	A	115	ILE	N-CA-C	-5.45	107.05	113.42
2	B	509	ARG	N-CA-C	-5.44	105.04	110.97
2	A	348	TYR	CA-C-N	5.36	124.54	118.97
2	A	348	TYR	C-N-CA	5.36	124.54	118.97
2	A	282	VAL	N-CA-C	5.25	115.42	107.80
2	B	302	GLU	N-CA-C	-5.23	101.25	109.25
2	B	443	GLY	N-CA-C	-5.20	100.86	113.18
2	A	64	GLY	N-CA-C	5.18	121.94	115.21
1	D	909	G	C2'-C3'-O3'	-5.13	106.00	113.70
2	A	272	LYS	N-CA-C	5.03	116.77	111.28

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	D	1517	0	768	36	0
2	A	4319	0	4339	357	0
2	B	3761	0	3774	263	0
3	A	22	0	10	5	0
All	All	9619	0	8891	634	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 34.

All (634) 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)
2:B:464:TYR:HA	2:B:500:ILE:HD11	1.32	1.07
2:B:448:LYS:H	2:B:448:LYS:HD3	1.20	1.05
2:B:55:ARG:HG2	2:B:55:ARG:HH21	1.34	0.92
2:B:47:HIS:HD2	2:B:49:LEU:H	1.13	0.92
2:A:460:THR:HG23	2:A:462:ILE:H	1.36	0.89
1:D:957:A:H1'	1:D:959:U:H5	1.37	0.88
2:A:484:GLN:HG2	2:A:485:GLU:H	1.41	0.86
2:A:200:ASP:HA	2:B:507:ASN:ND2	1.94	0.82
2:A:148:LEU:HG	2:A:160:ALA:HB2	1.63	0.81
1:D:909:G:H5'	1:D:910:G:OP2	1.80	0.80
2:B:385:VAL:HG13	2:B:465:ILE:HD12	1.61	0.80
2:A:106:GLY:HA2	2:A:110:GLU:HB2	1.63	0.80
2:A:380:ILE:O	2:A:384:ILE:HG13	1.82	0.79
2:B:47:HIS:CD2	2:B:49:LEU:H	1.98	0.79
2:B:286:HIS:HD2	2:B:288:LYS:H	1.30	0.79
2:A:227:TYR:HB3	2:B:73:ILE:HD12	1.66	0.78
1:D:957:A:H1'	1:D:959:U:C5	2.18	0.77
2:A:434:VAL:O	2:A:460:THR:HG22	1.84	0.77
2:B:55:ARG:HH21	2:B:55:ARG:CG	1.96	0.77
2:B:226:LEU:HD21	2:B:344:ARG:HG2	1.64	0.77
1:D:942:C:O2'	1:D:943:U:H5'	1.84	0.76
2:A:206:ILE:HB	2:A:237:VAL:HB	1.66	0.76
2:A:53:ILE:HG23	2:A:211:ILE:HG21	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:465:ILE:HD12	2:A:466:ASP:N	2.01	0.75
2:B:505:HIS:O	2:B:508:VAL:HG22	1.86	0.75
2:A:77:VAL:O	2:A:81:LYS:HG3	1.87	0.75
2:A:434:VAL:HB	2:A:460:THR:HG21	1.67	0.75
2:B:381:ALA:HA	2:B:384:ILE:HD12	1.69	0.75
2:A:374:THR:HG22	2:A:376:VAL:H	1.51	0.74
2:A:505:HIS:O	2:A:508:VAL:HG22	1.88	0.74
2:A:391:HIS:O	2:A:425:LEU:HD23	1.88	0.73
2:A:178:SER:HB3	2:B:176:PRO:HB2	1.71	0.72
2:B:490:LYS:HB3	2:B:528:THR:HG23	1.71	0.72
3:A:1002:SEP:N	3:A:1002:SEP:O2P	2.23	0.72
2:B:376:VAL:HA	2:B:379:LYS:HE2	1.71	0.72
2:B:267:ASP:HB3	2:B:279:GLN:HG2	1.70	0.71
2:B:399:SER:HA	2:B:417:VAL:HG12	1.71	0.71
2:B:504:ILE:HB	2:B:508:VAL:HG21	1.72	0.71
2:A:115:ILE:HD13	2:A:115:ILE:O	1.89	0.71
2:A:128:LEU:HD23	2:A:128:LEU:H	1.56	0.71
1:D:936:G:H21	2:A:522:ARG:H	1.38	0.71
2:A:37:PRO:O	2:A:364:ARG:HD3	1.90	0.71
2:B:448:LYS:HD3	2:B:448:LYS:N	2.03	0.71
2:B:381:ALA:HB2	2:B:473:ALA:HB2	1.73	0.71
2:B:504:ILE:HD12	2:B:508:VAL:HG23	1.74	0.70
2:A:97:LEU:HD23	2:B:95:PHE:CE2	2.27	0.69
2:B:263:ARG:HG2	2:B:300:TRP:CZ3	2.27	0.69
2:A:504:ILE:HB	2:A:508:VAL:HG21	1.76	0.68
1:D:934:C:H4'	1:D:935:A:OP2	1.90	0.68
2:A:151:GLU:HG3	2:A:152:VAL:HG13	1.76	0.68
1:D:959:U:H5'	1:D:960:C:OP2	1.93	0.68
2:A:118:ILE:HD12	2:A:118:ILE:H	1.57	0.68
2:A:464:TYR:HD1	2:A:464:TYR:H	1.42	0.67
2:A:326:LEU:HD13	2:A:327:GLY:N	2.09	0.67
2:A:475:LYS:HE2	2:A:475:LYS:HA	1.75	0.67
2:A:97:LEU:HD11	2:A:183:LEU:HG	1.76	0.67
2:B:53:ILE:HG23	2:B:211:ILE:HG21	1.75	0.67
2:A:219:GLN:O	2:A:221:GLU:HG3	1.95	0.67
2:A:434:VAL:HB	2:A:460:THR:CG2	2.25	0.67
2:A:31:SER:O	2:A:35:LEU:HD13	1.95	0.67
2:A:430:TYR:HD2	2:A:465:ILE:HG21	1.60	0.67
2:B:488:LYS:NZ	2:B:530:LYS:HE2	2.10	0.66
2:A:85:ARG:N	2:A:85:ARG:HD2	2.11	0.66
2:A:404:GLU:OE2	2:A:413:ARG:HB3	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:222:ASP:OD2	2:B:344:ARG:HD3	1.96	0.66
2:B:497:LEU:O	2:B:500:ILE:HG22	1.95	0.66
2:B:325:ASN:HD22	2:B:326:LEU:N	1.94	0.66
2:A:325:ASN:HD22	2:A:326:LEU:N	1.94	0.66
2:A:188:THR:HG23	2:A:309:TYR:OH	1.95	0.66
2:B:55:ARG:HG2	2:B:55:ARG:NH2	2.05	0.66
2:A:22:ALA:O	2:A:25:GLU:HG2	1.96	0.66
2:A:488:LYS:HD3	2:A:530:LYS:HE2	1.77	0.66
2:A:287:PRO:O	2:A:290:VAL:HG22	1.96	0.65
2:A:306:PHE:HB3	2:A:326:LEU:HD22	1.77	0.65
2:A:490:LYS:HD3	2:A:492:ARG:HH12	1.61	0.65
2:B:6:GLN:HG2	2:B:9:ARG:HH11	1.59	0.65
1:D:917:G:H4'	1:D:918:G:OP1	1.96	0.65
2:B:381:ALA:HB2	2:B:473:ALA:CB	2.27	0.65
2:A:85:ARG:HD2	2:A:85:ARG:H	1.61	0.65
2:A:385:VAL:HG13	2:A:465:ILE:HD13	1.79	0.65
2:B:263:ARG:HG2	2:B:300:TRP:CH2	2.32	0.65
2:B:465:ILE:HG13	2:B:466:ASP:H	1.62	0.64
2:B:437:TYR:CE2	2:B:438:LYS:HG2	2.32	0.64
2:A:139:GLU:HB3	2:A:143:ASP:OD1	1.97	0.64
2:A:387:THR:HG21	2:A:402:ALA:HA	1.80	0.64
2:B:216:ARG:HG2	2:B:218:GLU:HG2	1.80	0.64
2:A:397:PRO:O	2:A:398:CYS:HB3	1.99	0.63
2:B:360:LEU:HA	2:B:458:VAL:HG21	1.81	0.63
2:A:90:VAL:O	2:A:93:ARG:HG2	1.99	0.63
2:B:47:HIS:HD2	2:B:49:LEU:N	1.93	0.63
2:B:441:ILE:HB	2:B:519:ILE:HG23	1.80	0.63
2:A:414:VAL:HG13	2:A:531:ALA:HB2	1.81	0.63
2:A:137:LYS:HE3	2:A:137:LYS:HA	1.80	0.62
2:A:31:SER:HB2	2:A:32:PRO:HD2	1.81	0.62
2:B:37:PRO:O	2:B:364:ARG:HG3	1.99	0.62
2:A:178:SER:CB	2:B:176:PRO:HB2	2.28	0.62
2:B:485:GLU:HA	2:B:533:ILE:HG12	1.82	0.62
1:D:917:G:H2'	1:D:956:G:N2	2.14	0.62
2:B:494:VAL:O	2:B:522:ARG:HA	1.99	0.62
2:A:148:LEU:HG	2:A:160:ALA:CB	2.29	0.62
2:B:437:TYR:CD2	2:B:438:LYS:HG2	2.35	0.62
2:A:121:ARG:O	2:A:123:VAL:HG13	2.00	0.62
2:B:290:VAL:HA	2:B:297:SER:O	2.00	0.61
1:D:953:U:H3	1:D:957:A:H62	1.48	0.61
2:B:287:PRO:O	2:B:290:VAL:HG22	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:144:ASP:HA	2:A:147:TYR:CZ	2.36	0.61
2:A:397:PRO:CG	2:A:421:GLU:HG2	2.30	0.61
2:B:67:GLU:HG3	2:B:209:PHE:CZ	2.35	0.60
2:B:295:LYS:C	2:B:297:SER:H	2.07	0.60
2:A:485:GLU:HA	2:A:533:ILE:HG13	1.84	0.60
2:B:325:ASN:C	2:B:325:ASN:ND2	2.59	0.60
2:A:56:LEU:HD13	2:A:257:PHE:CZ	2.36	0.60
2:A:412:VAL:HG13	2:A:533:ILE:HG22	1.83	0.60
2:B:347:VAL:HG22	2:B:348:TYR:CD2	2.35	0.60
2:B:356:LYS:O	2:B:357:LEU:HD23	2.02	0.59
2:A:49:LEU:O	2:A:53:ILE:HG13	2.03	0.59
2:A:395:PRO:HA	2:A:423:THR:O	2.02	0.59
2:A:144:ASP:HA	2:A:147:TYR:CE1	2.38	0.59
2:B:372:PRO:HA	2:B:474:ARG:HD3	1.85	0.59
2:B:391:HIS:O	2:B:425:LEU:HD23	2.02	0.59
2:A:494:VAL:O	2:A:522:ARG:HA	2.03	0.59
2:A:504:ILE:HD12	2:A:508:VAL:HG23	1.84	0.59
2:B:325:ASN:HD22	2:B:325:ASN:C	2.11	0.58
2:B:11:LEU:O	2:B:11:LEU:HD23	2.03	0.58
2:B:432:ASN:HD21	2:B:445:PRO:HD3	1.68	0.58
2:B:451:ARG:O	2:B:454:PHE:N	2.37	0.58
2:B:508:VAL:O	2:B:511:TYR:HB3	2.04	0.58
2:A:3:PHE:HD2	2:A:26:ILE:HG21	1.68	0.58
2:A:213:ARG:HH21	2:A:228:THR:CG2	2.17	0.58
2:A:397:PRO:HD3	2:A:421:GLU:HG2	1.86	0.58
2:A:31:SER:CB	2:A:32:PRO:HD2	2.34	0.58
2:B:471:TYR:O	2:B:475:LYS:HB2	2.04	0.58
2:B:23:GLY:O	2:B:26:ILE:HG12	2.03	0.57
2:B:407:MET:HE1	2:B:477:GLU:HA	1.85	0.57
2:A:500:ILE:HG21	2:A:502:LEU:HD22	1.87	0.57
2:B:185:SER:O	2:B:214:CYS:HB3	2.04	0.57
2:A:464:TYR:CB	2:A:500:ILE:HD11	2.34	0.57
2:B:396:SER:OG	2:B:418:ASN:HB3	2.03	0.57
2:B:465:ILE:HG13	2:B:466:ASP:N	2.19	0.57
2:A:222:ASP:OD2	2:A:344:ARG:HG2	2.05	0.57
2:A:285:PHE:HB2	2:A:300:TRP:CE2	2.40	0.57
2:B:301:ILE:HD13	2:B:336:ILE:HD11	1.86	0.57
2:B:346:MET:O	2:B:349:PRO:HD3	2.05	0.57
2:A:314:LEU:HD11	2:A:323:VAL:HG12	1.87	0.57
2:A:397:PRO:HG3	2:A:421:GLU:HG2	1.87	0.56
2:A:213:ARG:HH21	2:A:228:THR:HG22	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:280:THR:HG22	2:B:281:GLU:N	2.20	0.56
2:A:184:ARG:NH1	2:A:212:ASP:HB3	2.20	0.56
2:A:267:ASP:HB2	2:A:281:GLU:HG3	1.86	0.56
2:B:174:LEU:C	2:B:175:LYS:HD2	2.30	0.56
2:A:311:PRO:HA	2:A:314:LEU:HD23	1.88	0.56
2:A:479:ALA:C	2:A:481:MET:H	2.14	0.56
2:B:7:LYS:O	2:B:11:LEU:HB2	2.05	0.56
2:A:115:ILE:HG23	2:A:116:GLU:HG3	1.88	0.56
2:B:346:MET:HE2	2:B:347:VAL:HG12	1.87	0.56
2:B:36:TYR:CD1	2:B:37:PRO:HA	2.40	0.56
2:B:388:ALA:O	2:B:392:ALA:HB2	2.05	0.56
2:A:306:PHE:CD1	2:A:306:PHE:C	2.83	0.56
2:B:356:LYS:HD2	2:B:357:LEU:H	1.71	0.56
2:B:412:VAL:HG22	2:B:533:ILE:HG22	1.87	0.56
2:A:309:TYR:HD1	2:A:323:VAL:CG1	2.18	0.55
2:A:320:PRO:HG2	2:A:321:TYR:CD1	2.41	0.55
2:A:154:ASP:HA	2:A:158:ILE:HD11	1.87	0.55
2:A:331:GLU:O	2:A:335:MET:HG3	2.06	0.55
2:B:175:LYS:HD2	2:B:175:LYS:N	2.21	0.55
1:D:912:A:H2'	1:D:913:G:O4'	2.06	0.55
2:A:325:ASN:ND2	2:A:325:ASN:C	2.60	0.55
2:B:85:ARG:HD2	2:B:85:ARG:N	2.22	0.55
2:B:102:LYS:NZ	2:B:102:LYS:HB3	2.21	0.55
2:A:430:TYR:HB2	2:A:465:ILE:CG2	2.37	0.55
2:A:504:ILE:O	2:A:504:ILE:HG13	2.07	0.55
2:B:101:PRO:O	2:B:103:PRO:HD3	2.07	0.55
2:A:156:ASP:CG	2:A:157:ASP:H	2.12	0.55
2:A:510:ARG:HH21	2:B:200:ASP:HB2	1.71	0.55
2:B:335:MET:SD	2:B:343:VAL:HG22	2.46	0.55
2:A:200:ASP:HA	2:B:507:ASN:HD22	1.71	0.55
2:A:243:VAL:HG12	2:A:247:LYS:HE3	1.88	0.55
2:A:325:ASN:HD22	2:A:325:ASN:C	2.14	0.55
2:A:492:ARG:HH11	2:A:492:ARG:HG3	1.71	0.55
2:B:451:ARG:O	2:B:452:SER:C	2.48	0.55
2:A:326:LEU:HD13	2:A:326:LEU:C	2.32	0.55
2:A:160:ALA:O	2:A:164:LEU:HG	2.07	0.55
2:A:458:VAL:HG23	2:A:458:VAL:O	2.07	0.54
2:B:376:VAL:HA	2:B:379:LYS:CE	2.36	0.54
2:A:173:GLU:O	2:A:173:GLU:HG3	2.08	0.54
2:B:432:ASN:ND2	2:B:445:PRO:HD3	2.23	0.54
2:A:202:LEU:HD21	2:B:355:ILE:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:183:LEU:HD11	2:B:183:LEU:HD11	1.90	0.54
2:A:484:GLN:NE2	2:A:487:VAL:HG22	2.22	0.54
2:B:206:ILE:HB	2:B:237:VAL:HB	1.90	0.54
2:A:258:GLY:HA3	2:A:286:HIS:CE1	2.42	0.54
2:A:344:ARG:HB3	2:A:352:HIS:CE1	2.42	0.54
2:B:376:VAL:O	2:B:380:ILE:HG13	2.07	0.54
2:A:200:ASP:HB2	2:B:510:ARG:HE	1.72	0.54
2:B:306:PHE:HB3	2:B:326:LEU:HD23	1.90	0.54
2:B:488:LYS:HZ2	2:B:530:LYS:HE2	1.72	0.54
2:A:135:TYR:HA	2:A:139:GLU:CG	2.38	0.54
2:A:478:GLU:HA	2:A:481:MET:HE3	1.90	0.54
2:A:90:VAL:O	2:A:92:ASP:N	2.41	0.54
2:A:347:VAL:HG22	2:A:348:TYR:CD2	2.42	0.54
2:B:218:GLU:C	2:B:220:GLY:H	2.16	0.54
2:A:236:LEU:HB3	2:A:324:MET:HB3	1.89	0.54
1:D:934:C:H5'	1:D:935:A:O5'	2.08	0.53
2:A:184:ARG:HH12	2:A:212:ASP:HB3	1.71	0.53
2:A:30:ARG:HB2	2:A:35:LEU:HD11	1.90	0.53
2:A:97:LEU:HD23	2:B:95:PHE:CD2	2.42	0.53
2:A:432:ASN:HB2	2:A:464:TYR:CE1	2.43	0.53
2:A:497:LEU:O	2:A:500:ILE:HG22	2.08	0.53
2:B:444:ILE:HG13	2:B:454:PHE:HE1	1.73	0.53
2:B:437:TYR:CD1	2:B:453:PHE:HA	2.43	0.53
2:B:496:ASN:ND2	2:B:499:ASP:H	2.05	0.53
1:D:917:G:HO2'	1:D:918:G:C5'	2.21	0.53
2:B:451:ARG:HH21	2:B:451:ARG:HG3	1.72	0.53
2:A:414:VAL:HG22	2:A:531:ALA:HB1	1.91	0.53
2:B:75:GLU:OE2	2:B:180:THR:HG23	2.08	0.53
2:B:264:PHE:HB3	2:B:280:THR:HG21	1.90	0.53
2:B:274:TYR:O	2:B:275:ILE:C	2.51	0.53
2:A:448:LYS:HD3	2:A:448:LYS:N	2.24	0.53
2:B:381:ALA:HA	2:B:384:ILE:CD1	2.38	0.53
2:A:225:ARG:HG3	2:A:226:LEU:H	1.74	0.53
2:A:442:TYR:CE2	2:A:453:PHE:HE2	2.27	0.53
2:A:259:PHE:CD2	2:A:286:HIS:HB2	2.44	0.52
2:A:464:TYR:HB3	2:A:500:ILE:HD11	1.92	0.52
2:B:197:HIS:O	2:B:201:LYS:HE3	2.08	0.52
2:A:264:PHE:HB3	2:A:280:THR:CG2	2.38	0.52
2:A:444:ILE:HB	2:A:450:TRP:CD1	2.44	0.52
1:D:916:G:H2'	1:D:959:U:O2'	2.09	0.52
1:D:925:G:H2'	1:D:926:G:O4'	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:30:ARG:HB2	2:A:35:LEU:CD1	2.39	0.52
2:B:448:LYS:H	2:B:448:LYS:CD	2.01	0.52
2:A:218:GLU:HG3	2:A:225:ARG:HG2	1.91	0.52
2:A:188:THR:HG23	2:A:309:TYR:CZ	2.43	0.52
2:A:333:LEU:HD22	2:A:337:LEU:HD11	1.92	0.52
2:A:401:LEU:HD13	2:A:415:TYR:HE1	1.74	0.52
2:B:416:VAL:HG22	2:B:529:VAL:HG22	1.90	0.52
2:B:479:ALA:HB1	2:B:484:GLN:HB2	1.92	0.52
1:D:925:G:H2'	1:D:926:G:H8	1.75	0.52
2:A:118:ILE:H	2:A:118:ILE:CD1	2.22	0.52
2:B:320:PRO:HG2	2:B:321:TYR:CD1	2.45	0.52
2:B:331:GLU:O	2:B:335:MET:HG3	2.10	0.52
2:B:364:ARG:HG3	2:B:364:ARG:HH21	1.74	0.52
2:B:458:VAL:HG23	2:B:458:VAL:O	2.08	0.52
2:A:407:MET:HG2	2:A:408:MET:SD	2.50	0.52
2:A:430:TYR:CD2	2:A:465:ILE:HG21	2.42	0.52
2:B:392:ALA:HB1	2:B:426:CYS:HB2	1.91	0.52
2:A:327:GLY:HA3	3:A:1001:SEP:HB2	1.92	0.51
2:B:53:ILE:HG23	2:B:211:ILE:CG2	2.40	0.51
2:B:363:ALA:HB2	2:B:458:VAL:HG23	1.92	0.51
2:A:408:MET:HE2	2:A:480:ALA:HB3	1.92	0.51
2:B:482:ARG:O	2:B:483:GLU:HB2	2.11	0.51
2:A:394:GLU:HG3	2:A:400:PHE:CZ	2.46	0.51
2:B:430:TYR:HD2	2:B:465:ILE:HG21	1.74	0.51
2:B:504:ILE:HB	2:B:508:VAL:CG2	2.39	0.51
2:B:396:SER:HB3	2:B:418:ASN:O	2.09	0.51
2:A:185:SER:HA	2:A:215:PHE:O	2.11	0.51
2:A:287:PRO:O	2:A:289:LEU:N	2.43	0.51
2:A:413:ARG:NH1	2:A:532:GLU:HG3	2.24	0.51
1:D:925:G:H2'	1:D:926:G:C8	2.45	0.51
2:A:118:ILE:HD12	2:A:118:ILE:N	2.25	0.51
2:A:432:ASN:ND2	2:A:445:PRO:HD3	2.26	0.51
2:B:38:ARG:HG2	2:B:367:LYS:HG2	1.91	0.51
2:B:444:ILE:HG13	2:B:454:PHE:CE1	2.45	0.51
2:A:306:PHE:HB3	2:A:326:LEU:CD2	2.39	0.51
2:A:92:ASP:O	2:A:217:ARG:HD3	2.12	0.51
2:B:412:VAL:CG2	2:B:533:ILE:HG22	2.41	0.50
2:B:484:GLN:O	2:B:533:ILE:HD11	2.11	0.50
2:A:309:TYR:CD1	2:A:323:VAL:HG13	2.47	0.50
2:A:397:PRO:CD	2:A:421:GLU:HG2	2.41	0.50
2:B:347:VAL:C	2:B:349:PRO:HD2	2.36	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:916:G:O5'	1:D:916:G:H8	1.95	0.50
2:A:238:ASP:O	2:A:322:PRO:HD2	2.12	0.50
2:B:261:ASN:HD22	2:B:262:PHE:H	1.59	0.50
2:B:496:ASN:C	2:B:496:ASN:HD22	2.18	0.50
2:A:198:ILE:HD12	2:A:202:LEU:HD12	1.93	0.50
2:A:406:GLU:OE1	2:A:409:GLY:HA2	2.12	0.50
2:B:72:LEU:HA	2:B:184:ARG:NH2	2.26	0.50
2:B:83:PHE:CD2	2:B:90:VAL:HG21	2.46	0.50
2:B:478:GLU:O	2:B:482:ARG:HB2	2.11	0.50
2:A:158:ILE:HG13	2:A:159:THR:HG23	1.93	0.50
2:A:161:VAL:HG23	2:A:162:LYS:N	2.27	0.50
2:B:213:ARG:HH21	2:B:228:THR:HG22	1.76	0.50
2:A:47:HIS:CD2	2:A:49:LEU:HB2	2.46	0.50
2:A:53:ILE:HG23	2:A:211:ILE:CG2	2.39	0.50
2:B:185:SER:HA	2:B:215:PHE:O	2.11	0.50
2:A:16:PHE:O	2:A:19:ALA:N	2.44	0.50
2:A:478:GLU:HB3	2:A:482:ARG:NH2	2.26	0.50
2:A:496:ASN:C	2:A:496:ASN:HD22	2.20	0.50
2:B:450:TRP:O	2:B:451:ARG:C	2.54	0.50
2:A:347:VAL:O	2:A:349:PRO:HD2	2.12	0.50
2:B:340:TYR:CE2	2:B:346:MET:HA	2.46	0.50
2:A:294:THR:HG23	2:A:296:TYR:H	1.77	0.49
2:A:410:ARG:O	2:A:412:VAL:HG23	2.12	0.49
2:B:391:HIS:C	2:B:393:SER:H	2.20	0.49
2:A:23:GLY:O	2:A:25:GLU:N	2.45	0.49
2:A:40:GLY:HA3	2:A:364:ARG:NH2	2.26	0.49
2:B:290:VAL:HG23	2:B:290:VAL:O	2.13	0.49
1:D:915:G:H2'	1:D:916:G:O4'	2.13	0.49
2:B:47:HIS:CD2	2:B:49:LEU:HB2	2.48	0.49
2:B:437:TYR:HD2	2:B:438:LYS:N	2.11	0.49
1:D:923:G:O2'	1:D:924:C:H5'	2.13	0.49
2:A:401:LEU:HA	2:A:415:TYR:HD1	1.77	0.49
2:A:435:VAL:HG21	2:A:444:ILE:HD11	1.95	0.49
2:A:227:TYR:HB3	2:B:73:ILE:CD1	2.41	0.49
2:A:510:ARG:HE	2:B:200:ASP:HB2	1.78	0.49
2:B:363:ALA:HB2	2:B:458:VAL:CG2	2.42	0.49
1:D:967:C:H2'	1:D:968:U:O4'	2.13	0.49
2:A:188:THR:HA	2:A:191:TRP:HD1	1.78	0.49
2:A:381:ALA:HB2	2:A:473:ALA:HB2	1.95	0.49
2:B:309:TYR:HB2	2:B:314:LEU:HD21	1.93	0.49
2:B:321:TYR:HB3	2:B:322:PRO:HD2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:519:ILE:HG22	2:B:521:VAL:HG23	1.95	0.48
2:A:460:THR:HG23	2:A:462:ILE:N	2.17	0.48
2:B:209:PHE:HA	2:B:234:CYS:HA	1.96	0.48
2:B:222:ASP:OD2	2:B:226:LEU:HD23	2.14	0.48
2:A:283:PHE:CE1	2:A:302:GLU:HB2	2.49	0.48
2:A:324:MET:HE1	2:A:326:LEU:HB2	1.96	0.48
2:A:372:PRO:HG3	2:A:473:ALA:HB3	1.96	0.48
2:A:525:LEU:HD22	2:A:527:VAL:HG12	1.94	0.48
2:B:246:GLY:HA3	2:B:306:PHE:CE1	2.49	0.48
2:A:155:VAL:HG22	2:A:158:ILE:HD13	1.96	0.48
2:A:258:GLY:O	2:A:259:PHE:C	2.57	0.48
2:A:271:SER:HB3	2:A:273:TYR:CE1	2.48	0.48
2:A:463:ARG:HB2	2:A:466:ASP:OD2	2.13	0.48
2:A:68:VAL:HG13	2:A:69:VAL:N	2.27	0.48
2:A:154:ASP:CA	2:A:158:ILE:HD11	2.43	0.48
2:B:376:VAL:HG12	2:B:380:ILE:HD11	1.95	0.48
2:A:285:PHE:HB2	2:A:300:TRP:CD2	2.49	0.48
2:A:351:ILE:HG22	2:A:352:HIS:HD2	1.79	0.48
2:A:478:GLU:HB3	2:A:482:ARG:HH22	1.77	0.48
2:A:264:PHE:HB3	2:A:280:THR:HG21	1.94	0.48
2:A:315:ALA:C	2:A:317:TYR:N	2.71	0.48
2:A:504:ILE:HB	2:A:508:VAL:CG2	2.42	0.48
2:A:187:MET:HG2	2:A:233:SER:HB3	1.95	0.48
2:A:222:ASP:CG	2:A:344:ARG:HG2	2.39	0.48
2:A:447:THR:C	2:A:449:LYS:H	2.20	0.48
2:B:344:ARG:HB3	2:B:352:HIS:NE2	2.28	0.48
2:B:433:GLU:OE2	2:B:459:PRO:HB2	2.13	0.48
2:A:127:PRO:C	2:A:129:GLN:H	2.21	0.47
2:A:267:ASP:HB3	2:A:279:GLN:HG2	1.96	0.47
2:A:148:LEU:HG	2:A:160:ALA:CA	2.44	0.47
2:A:429:ALA:HA	2:A:432:ASN:OD1	2.14	0.47
2:A:498:SER:C	2:A:500:ILE:H	2.22	0.47
2:B:19:ALA:O	2:B:22:ALA:HB3	2.14	0.47
2:B:254:LEU:O	2:B:259:PHE:HD1	1.97	0.47
2:A:47:HIS:CD2	2:A:49:LEU:H	2.33	0.47
2:B:372:PRO:HB2	2:B:378:LEU:HD23	1.95	0.47
2:A:52:THR:HG23	2:A:337:LEU:HD12	1.97	0.47
2:A:369:LYS:N	2:A:501:ASN:O	2.44	0.47
3:A:1001:SEP:P	3:A:1001:SEP:H	2.37	0.47
2:A:97:LEU:N	2:A:97:LEU:HD12	2.29	0.47
2:A:380:ILE:HG22	2:A:384:ILE:HD11	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:267:ASP:CB	2:B:279:GLN:HG2	2.43	0.47
2:B:421:GLU:O	2:B:422:ASN:C	2.57	0.47
2:A:309:TYR:HD1	2:A:323:VAL:HG13	1.79	0.47
2:A:464:TYR:CD1	2:A:464:TYR:N	2.81	0.47
2:B:213:ARG:HH21	2:B:228:THR:CG2	2.28	0.47
2:B:237:VAL:HG22	2:B:323:VAL:HG22	1.97	0.47
2:A:93:ARG:O	2:A:217:ARG:HG2	2.15	0.47
2:A:214:CYS:C	2:A:215:PHE:CD1	2.92	0.47
2:A:298:ASP:OD2	2:A:298:ASP:N	2.48	0.47
2:A:315:ALA:C	2:A:317:TYR:H	2.23	0.47
2:A:479:ALA:C	2:A:481:MET:N	2.71	0.47
2:B:324:MET:SD	2:B:324:MET:C	2.97	0.47
2:B:328:LEU:HD12	2:B:328:LEU:C	2.40	0.47
2:B:330:VAL:O	2:B:331:GLU:C	2.57	0.47
2:B:436:VAL:HG12	2:B:437:TYR:N	2.29	0.47
2:B:490:LYS:HA	2:B:528:THR:HA	1.96	0.47
2:A:149:ILE:HA	2:A:154:ASP:O	2.15	0.47
2:A:489:VAL:O	2:A:528:THR:HG23	2.15	0.47
2:B:6:GLN:O	2:B:7:LYS:C	2.56	0.47
2:B:32:PRO:HD2	2:B:371:VAL:HG11	1.96	0.47
2:B:275:ILE:HG13	2:B:310:SER:HA	1.95	0.47
2:B:314:LEU:HD12	2:B:319:ILE:HG21	1.96	0.47
2:A:112:ILE:O	2:A:112:ILE:HG13	2.15	0.47
2:A:401:LEU:HD13	2:A:415:TYR:CE1	2.50	0.47
2:B:383:SER:O	2:B:384:ILE:C	2.58	0.47
2:A:333:LEU:O	2:A:337:LEU:HG	2.14	0.47
2:A:102:LYS:HD2	2:A:135:TYR:CE1	2.50	0.46
2:B:32:PRO:O	2:B:38:ARG:NH1	2.48	0.46
2:B:295:LYS:C	2:B:297:SER:N	2.73	0.46
2:B:306:PHE:HB3	2:B:326:LEU:CD2	2.45	0.46
2:B:385:VAL:O	2:B:389:GLU:HG3	2.15	0.46
2:A:430:TYR:O	2:A:465:ILE:HG23	2.14	0.46
1:D:917:G:O2'	1:D:918:G:O5'	2.30	0.46
2:A:331:GLU:H	2:A:331:GLU:CD	2.24	0.46
2:A:81:LYS:HE3	2:A:316:GLU:HB3	1.98	0.46
2:B:269:LYS:O	2:B:270:ARG:C	2.58	0.46
2:B:405:GLY:O	2:B:412:VAL:N	2.49	0.46
2:A:285:PHE:CD1	2:A:285:PHE:C	2.93	0.46
2:A:310:SER:C	2:A:312:THR:N	2.73	0.46
2:A:310:SER:C	2:A:312:THR:H	2.24	0.46
2:A:450:TRP:O	2:A:451:ARG:C	2.59	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:500:ILE:HG23	2:A:502:LEU:HD13	1.98	0.46
2:B:388:ALA:HB1	2:B:426:CYS:SG	2.56	0.46
2:A:306:PHE:HB2	2:A:325:ASN:O	2.16	0.46
2:A:463:ARG:O	2:A:464:TYR:C	2.59	0.46
2:A:466:ASP:O	2:A:470:TYR:CD2	2.69	0.46
2:B:318:ASP:O	2:B:320:PRO:HD3	2.15	0.46
2:A:72:LEU:C	2:A:72:LEU:HD23	2.41	0.46
2:A:294:THR:C	2:A:296:TYR:H	2.22	0.46
2:A:306:PHE:HB2	2:A:324:MET:CE	2.45	0.46
2:B:242:SER:C	2:B:244:ASP:N	2.73	0.46
2:B:342:ASP:O	2:B:343:VAL:C	2.59	0.46
2:B:410:ARG:HD3	2:B:410:ARG:N	2.30	0.46
2:A:131:ILE:HD12	2:A:131:ILE:N	2.31	0.45
2:A:135:TYR:HA	2:A:139:GLU:HG3	1.98	0.45
2:A:401:LEU:HA	2:A:415:TYR:CD1	2.52	0.45
2:A:533:ILE:HG13	2:A:533:ILE:O	2.16	0.45
2:B:71:PRO:HG2	2:B:193:ILE:HG21	1.98	0.45
2:A:222:ASP:HA	2:A:225:ARG:O	2.16	0.45
2:A:258:GLY:HA3	2:A:286:HIS:HE1	1.79	0.45
2:A:484:GLN:HG2	2:A:485:GLU:N	2.20	0.45
2:B:188:THR:HA	2:B:191:TRP:HD1	1.80	0.45
2:A:287:PRO:C	2:A:289:LEU:N	2.74	0.45
2:A:125:SER:O	2:A:126:LYS:C	2.60	0.45
2:A:184:ARG:NH2	2:A:187:MET:SD	2.90	0.45
2:A:349:PRO:HG2	2:A:350:GLN:OE1	2.16	0.45
2:A:409:GLY:C	2:A:410:ARG:HG3	2.41	0.45
2:A:414:VAL:HG22	2:A:531:ALA:CB	2.47	0.45
2:B:99:THR:OG1	2:B:100:LEU:N	2.49	0.45
1:D:957:A:HO2'	1:D:958:A:P	2.40	0.45
2:A:434:VAL:HG21	2:A:502:LEU:CD2	2.47	0.45
1:D:931:C:O2'	1:D:932:U:H5'	2.17	0.45
2:A:95:PHE:CE2	2:B:97:LEU:HD23	2.51	0.45
2:A:504:ILE:O	2:A:505:HIS:C	2.59	0.45
2:B:354:GLU:O	2:B:355:ILE:C	2.59	0.45
2:A:145:LEU:HD22	2:A:145:LEU:N	2.32	0.45
2:A:217:ARG:NH2	2:B:99:THR:OG1	2.49	0.45
2:A:257:PHE:O	2:A:286:HIS:HE1	2.00	0.45
2:A:511:TYR:CZ	2:A:515:LYS:HG3	2.52	0.45
2:B:393:SER:O	2:B:394:GLU:C	2.60	0.45
2:B:442:TYR:CE2	2:B:453:PHE:HE2	2.35	0.45
1:D:941:G:C2'	1:D:942:C:H5'	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:957:A:O2'	1:D:958:A:P	2.75	0.45
2:A:399:SER:HA	2:A:417:VAL:HG12	1.99	0.44
2:A:377:GLY:O	2:A:380:ILE:HB	2.17	0.44
2:A:500:ILE:CG2	2:A:502:LEU:HB2	2.47	0.44
2:B:188:THR:HG23	2:B:309:TYR:CZ	2.52	0.44
2:B:275:ILE:HG12	2:B:311:PRO:HD3	1.99	0.44
2:A:434:VAL:HG21	2:A:502:LEU:HD23	1.99	0.44
2:B:75:GLU:HB3	2:B:77:VAL:HG12	1.99	0.44
2:A:350:GLN:H	2:A:350:GLN:CD	2.26	0.44
2:B:16:PHE:O	2:B:19:ALA:HB3	2.17	0.44
2:A:468:PHE:O	2:A:471:TYR:N	2.48	0.44
2:B:23:GLY:C	2:B:25:GLU:H	2.25	0.44
2:A:200:ASP:HB2	2:B:510:ARG:HH21	1.83	0.44
2:A:403:PHE:O	2:A:413:ARG:HA	2.17	0.44
2:B:251:GLU:HG3	2:B:262:PHE:CE2	2.51	0.44
2:B:527:VAL:O	2:B:527:VAL:HG13	2.17	0.44
2:B:282:VAL:CG2	2:B:303:ILE:HG22	2.47	0.44
2:B:407:MET:HE2	2:B:480:ALA:CB	2.48	0.44
2:A:67:GLU:HG3	2:A:209:PHE:CZ	2.52	0.44
2:A:176:PRO:HG2	2:B:178:SER:CB	2.48	0.44
2:A:186:HIS:HB2	3:A:1001:SEP:O2P	2.18	0.44
2:A:433:GLU:OE2	2:A:461:GLY:HA2	2.18	0.44
2:A:465:ILE:HD12	2:A:465:ILE:C	2.42	0.44
2:A:453:PHE:N	2:A:453:PHE:CD1	2.86	0.43
2:B:377:GLY:O	2:B:380:ILE:HB	2.18	0.43
2:B:326:LEU:C	2:B:326:LEU:HD13	2.43	0.43
2:B:485:GLU:HA	2:B:533:ILE:CG1	2.48	0.43
2:B:506:GLU:HB2	2:B:510:ARG:HH12	1.83	0.43
2:A:56:LEU:HB3	2:A:211:ILE:CD1	2.49	0.43
2:A:286:HIS:CD2	2:A:288:LYS:HB2	2.53	0.43
2:A:409:GLY:O	2:A:410:ARG:HG3	2.17	0.43
2:B:479:ALA:HA	2:B:482:ARG:HB3	2.00	0.43
2:A:85:ARG:C	2:A:87:ALA:H	2.26	0.43
2:A:118:ILE:HG22	2:A:119:THR:N	2.34	0.43
2:A:141:ASP:CG	2:A:142:GLY:N	2.77	0.43
2:A:215:PHE:CD1	2:A:215:PHE:N	2.86	0.43
2:A:251:GLU:HG3	2:A:262:PHE:CE2	2.53	0.43
2:A:273:TYR:O	2:A:309:TYR:HA	2.19	0.43
2:A:500:ILE:CG2	2:A:502:LEU:HD22	2.47	0.43
2:B:261:ASN:HD22	2:B:262:PHE:N	2.16	0.43
2:B:275:ILE:CG1	2:B:310:SER:HA	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:447:THR:C	2:B:449:LYS:N	2.75	0.43
2:B:472:ALA:O	2:B:475:LYS:N	2.45	0.43
2:B:525:LEU:HD13	2:B:527:VAL:HG12	2.01	0.43
2:A:153:LEU:N	2:A:153:LEU:HD12	2.34	0.43
2:A:155:VAL:O	2:A:156:ASP:C	2.61	0.43
2:A:162:LYS:O	2:A:166:GLU:HG2	2.19	0.43
2:A:328:LEU:HD11	2:A:333:LEU:HD12	2.00	0.43
2:A:472:ALA:O	2:A:475:LYS:N	2.51	0.43
2:B:340:TYR:CD2	2:B:346:MET:HA	2.53	0.43
2:B:368:VAL:HG11	2:B:371:VAL:HG22	2.01	0.43
2:A:6:GLN:O	2:A:7:LYS:C	2.61	0.43
2:A:135:TYR:HA	2:A:139:GLU:HG2	2.00	0.43
2:A:286:HIS:HD2	2:A:288:LYS:HB2	1.82	0.43
2:A:303:ILE:HG13	2:A:328:LEU:HD21	1.99	0.43
2:A:351:ILE:HG22	2:A:352:HIS:CD2	2.53	0.43
2:B:292:SER:OG	2:B:294:THR:HG22	2.19	0.43
1:D:957:A:O2'	1:D:958:A:O5'	2.28	0.43
2:A:124:ASP:CG	2:A:125:SER:H	2.26	0.43
2:A:157:ASP:HA	2:A:160:ALA:HB3	2.00	0.43
2:A:187:MET:CB	3:A:1001:SEP:O3P	2.67	0.43
2:A:515:LYS:O	2:A:516:LYS:HB2	2.18	0.43
2:B:23:GLY:C	2:B:25:GLU:N	2.77	0.43
2:B:294:THR:OG1	2:B:295:LYS:N	2.52	0.43
2:B:407:MET:C	2:B:409:GLY:H	2.26	0.43
2:A:90:VAL:C	2:A:92:ASP:H	2.27	0.43
2:A:145:LEU:HD12	2:A:156:ASP:HB2	2.00	0.43
2:A:52:THR:O	2:A:56:LEU:CD2	2.66	0.43
2:A:342:ASP:O	2:A:343:VAL:C	2.62	0.43
2:A:381:ALA:O	2:A:382:GLN:C	2.62	0.43
2:B:282:VAL:HG23	2:B:303:ILE:HG22	2.01	0.43
2:B:421:GLU:O	2:B:423:THR:HG22	2.18	0.43
2:B:451:ARG:HG3	2:B:451:ARG:NH2	2.34	0.43
1:D:968:U:H2'	1:D:969:G:C8	2.54	0.43
2:A:155:VAL:HG23	2:A:158:ILE:CG2	2.49	0.43
2:A:213:ARG:NH2	2:A:228:THR:CG2	2.81	0.43
2:A:243:VAL:HG13	2:A:306:PHE:CE1	2.54	0.43
2:B:475:LYS:HD2	2:B:478:GLU:OE2	2.18	0.43
1:D:943:U:H2'	1:D:944:U:H5'	2.00	0.42
2:A:8:TYR:HH	2:A:25:GLU:CD	2.26	0.42
2:A:16:PHE:O	2:A:17:GLU:C	2.60	0.42
2:A:251:GLU:HG2	2:A:255:ARG:HG3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:468:PHE:O	2:A:471:TYR:HB3	2.18	0.42
2:B:85:ARG:C	2:B:87:ALA:H	2.27	0.42
2:B:183:LEU:HD13	2:B:215:PHE:CD1	2.54	0.42
1:D:941:G:O2'	1:D:942:C:H5'	2.19	0.42
2:A:153:LEU:O	2:A:154:ASP:HB2	2.19	0.42
2:A:370:GLU:O	2:A:501:ASN:ND2	2.45	0.42
2:A:475:LYS:HE2	2:A:475:LYS:CA	2.48	0.42
2:A:397:PRO:HD3	2:A:421:GLU:CG	2.49	0.42
2:B:90:VAL:O	2:B:93:ARG:HG2	2.19	0.42
2:B:286:HIS:CD2	2:B:288:LYS:HB3	2.54	0.42
2:B:486:GLU:HA	2:B:531:ALA:O	2.19	0.42
1:D:936:G:H2'	2:A:522:ARG:CZ	2.50	0.42
2:A:464:TYR:HD1	2:A:464:TYR:N	2.13	0.42
2:A:137:LYS:HE3	2:A:137:LYS:CA	2.47	0.42
2:A:337:LEU:HB2	2:A:338:TYR:CD1	2.55	0.42
2:A:380:ILE:O	2:A:383:SER:OG	2.27	0.42
2:B:275:ILE:HG13	2:B:309:TYR:O	2.19	0.42
2:A:56:LEU:HD13	2:A:257:PHE:HZ	1.84	0.42
2:A:273:TYR:CE2	2:A:274:TYR:CE2	3.07	0.42
2:A:292:SER:OG	2:A:294:THR:HG22	2.19	0.42
2:A:321:TYR:HB3	2:A:322:PRO:HD2	2.02	0.42
2:A:325:ASN:ND2	2:A:326:LEU:N	2.63	0.42
2:B:437:TYR:HE2	2:B:438:LYS:NZ	2.17	0.42
2:A:239:GLU:HG2	2:A:321:TYR:CE2	2.53	0.42
2:A:500:ILE:HG22	2:A:502:LEU:H	1.85	0.42
2:B:23:GLY:O	2:B:25:GLU:N	2.52	0.42
2:B:445:PRO:O	2:B:447:THR:N	2.53	0.42
2:B:496:ASN:HD21	2:B:499:ASP:H	1.67	0.42
2:A:447:THR:C	2:A:449:LYS:N	2.77	0.42
2:A:488:LYS:HG2	2:A:530:LYS:HG2	2.01	0.42
2:B:354:GLU:OE2	2:B:354:GLU:HA	2.20	0.42
1:D:909:G:H1'	1:D:946:A:H1'	2.01	0.42
2:A:121:ARG:O	2:A:122:GLU:C	2.63	0.42
2:B:204:LEU:HD23	2:B:238:ASP:HA	2.02	0.42
1:D:968:U:H2'	1:D:969:G:H8	1.85	0.41
2:B:56:LEU:HB3	2:B:211:ILE:CD1	2.51	0.41
2:B:297:SER:C	2:B:299:GLY:N	2.78	0.41
2:A:90:VAL:C	2:A:92:ASP:N	2.77	0.41
2:A:155:VAL:HG23	2:A:158:ILE:HG23	2.02	0.41
2:A:156:ASP:CG	2:A:157:ASP:N	2.78	0.41
2:A:394:GLU:HG3	2:A:400:PHE:HZ	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:511:TYR:CE1	2:A:515:LYS:HG3	2.54	0.41
2:B:282:VAL:HG13	2:B:304:ALA:O	2.19	0.41
2:B:324:MET:SD	2:B:325:ASN:N	2.93	0.41
2:A:326:LEU:C	2:A:326:LEU:CD1	2.92	0.41
2:A:448:LYS:N	2:A:448:LYS:CD	2.83	0.41
2:B:271:SER:HB3	2:B:273:TYR:CE1	2.55	0.41
2:B:320:PRO:HG2	2:B:321:TYR:CE1	2.54	0.41
2:A:140:ILE:N	2:A:143:ASP:OD1	2.50	0.41
2:A:445:PRO:O	2:A:450:TRP:HD1	2.04	0.41
2:B:102:LYS:HB3	2:B:102:LYS:HZ2	1.84	0.41
2:A:56:LEU:HB3	2:A:211:ILE:HD11	2.03	0.41
2:A:484:GLN:HE22	2:A:487:VAL:HG22	1.83	0.41
2:A:507:ASN:ND2	2:B:200:ASP:C	2.78	0.41
2:B:47:HIS:CG	2:B:48:PRO:HD2	2.55	0.41
2:B:235:VAL:HA	2:B:324:MET:O	2.20	0.41
2:B:281:GLU:HG2	2:B:305:THR:HG22	2.02	0.41
1:D:917:G:O2'	1:D:918:G:C5'	2.68	0.41
2:A:6:GLN:O	2:A:10:GLU:HG2	2.21	0.41
2:A:203:PRO:HG3	2:B:508:VAL:HG12	2.02	0.41
2:A:366:ILE:HA	2:A:504:ILE:HG22	2.03	0.41
2:A:453:PHE:N	2:A:453:PHE:HD1	2.19	0.41
2:A:330:VAL:O	2:A:331:GLU:C	2.64	0.41
2:A:508:VAL:HG12	2:B:203:PRO:HG3	2.02	0.41
2:B:288:LYS:NZ	2:B:337:LEU:O	2.50	0.41
1:D:934:C:H6	1:D:934:C:O5'	2.03	0.41
2:A:3:PHE:CD1	2:A:3:PHE:C	2.98	0.41
2:A:95:PHE:CD2	2:B:97:LEU:HD23	2.56	0.41
2:A:101:PRO:HB3	2:A:177:ILE:CD1	2.51	0.41
2:A:309:TYR:HD1	2:A:323:VAL:HG11	1.82	0.41
2:B:10:GLU:HG3	2:B:11:LEU:N	2.36	0.41
2:B:360:LEU:O	2:B:363:ALA:HB3	2.21	0.41
2:B:423:THR:HG23	2:B:424:LYS:N	2.36	0.41
1:D:918:G:H5''	1:D:958:A:H61	1.85	0.41
2:A:16:PHE:O	2:A:18:ALA:N	2.54	0.41
2:A:38:ARG:NH2	2:A:460:THR:O	2.54	0.41
2:A:145:LEU:CD1	2:A:156:ASP:HB2	2.51	0.41
2:A:211:ILE:HG12	2:A:232:ALA:HA	2.02	0.41
2:A:346:MET:SD	2:A:346:MET:C	3.04	0.41
2:A:397:PRO:O	2:A:398:CYS:CB	2.68	0.41
2:A:520:ASP:OD1	2:A:522:ARG:NH2	2.54	0.41
2:B:77:VAL:HG22	2:B:81:LYS:HD2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:218:GLU:OE2	2:B:218:GLU:HA	2.21	0.41
2:B:381:ALA:O	2:B:384:ILE:HB	2.19	0.41
2:B:492:ARG:HA	2:B:526:PHE:HD2	1.86	0.41
2:A:235:VAL:HA	2:A:324:MET:O	2.21	0.41
2:A:383:SER:HA	2:A:386:GLU:HB3	2.03	0.41
2:A:500:ILE:HG23	2:A:500:ILE:O	2.21	0.41
2:B:1:MET:N	2:B:26:ILE:O	2.54	0.41
2:B:378:LEU:C	2:B:380:ILE:N	2.79	0.41
2:A:67:GLU:HB2	2:B:46:GLU:HA	2.03	0.40
2:A:259:PHE:CE2	2:A:286:HIS:HB2	2.57	0.40
2:A:314:LEU:HD11	2:A:323:VAL:CG1	2.49	0.40
2:B:32:PRO:HG3	2:B:470:TYR:CZ	2.56	0.40
2:B:373:GLN:OE1	2:B:373:GLN:HA	2.21	0.40
2:B:396:SER:HA	2:B:397:PRO:C	2.46	0.40
2:B:513:LEU:HD12	2:B:513:LEU:HA	1.92	0.40
2:A:121:ARG:HG3	2:A:121:ARG:HH11	1.86	0.40
2:A:356:LYS:O	2:A:357:LEU:HD23	2.20	0.40
2:A:468:PHE:O	2:A:469:ALA:C	2.62	0.40
2:B:65:PHE:CD1	2:B:65:PHE:N	2.89	0.40
2:B:92:ASP:O	2:B:217:ARG:HD3	2.21	0.40
2:B:188:THR:HG23	2:B:309:TYR:OH	2.21	0.40
2:B:268:GLU:O	2:B:270:ARG:HG3	2.20	0.40
2:A:80:LYS:HA	2:A:87:ALA:CB	2.52	0.40
2:A:106:GLY:CA	2:A:110:GLU:HB2	2.43	0.40
2:A:141:ASP:C	2:A:143:ASP:H	2.30	0.40
2:B:211:ILE:HA	2:B:231:SER:O	2.21	0.40
2:A:23:GLY:C	2:A:25:GLU:N	2.79	0.40
2:A:37:PRO:C	2:A:39:VAL:N	2.78	0.40
2:A:47:HIS:CG	2:A:48:PRO:HD2	2.56	0.40
2:A:85:ARG:N	2:A:85:ARG:CD	2.83	0.40
2:A:290:VAL:HA	2:A:297:SER:O	2.22	0.40
2:B:271:SER:HB3	2:B:273:TYR:CD1	2.57	0.40
2:B:430:TYR:HB2	2:B:465:ILE:HG23	2.03	0.40
2:A:277:ASP:O	2:A:277:ASP:CG	2.64	0.40
2:A:423:THR:OG1	2:A:424:LYS:N	2.54	0.40
2:B:387:THR:C	2:B:389:GLU:H	2.30	0.40
2:B:447:THR:O	2:B:449:LYS:N	2.54	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	
2	A	532/534 (100%)	417 (78%)	88 (16%)	27 (5%)	1	13
2	B	461/534 (86%)	376 (82%)	71 (15%)	14 (3%)	3	23
All	All	993/1068 (93%)	793 (80%)	159 (16%)	41 (4%)	2	17

All (41) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	118	ILE
2	A	398	CYS
2	A	526	PHE
2	B	14	LYS
2	B	484	GLN
2	A	24	LYS
2	A	91	LEU
2	A	107	ILE
2	A	125	SER
2	A	156	ASP
2	A	288	LYS
2	A	484	GLN
2	A	500	ILE
2	B	423	THR
2	A	148	LEU
2	A	151	GLU
2	A	199	ALA
2	B	24	LYS
2	B	448	LYS
2	B	451	ARG
2	B	526	PHE
2	A	16	PHE
2	A	17	GLU
2	A	104	ASN
2	A	122	GLU

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Mol	Chain	Res	Type
2	A	369	LYS
2	B	422	ASN
2	A	154	ASP
2	A	339	GLY
2	A	279	GLN
2	B	353	GLY
2	B	409	GLY
2	B	446	LYS
2	A	140	ILE
2	A	343	VAL
2	B	343	VAL
2	A	105	VAL
2	A	126	LYS
2	A	127	PRO
2	B	103	PRO
2	B	417	VAL

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	
2	A	464/464 (100%)	442 (95%)	22 (5%)	23	57
2	B	402/464 (87%)	375 (93%)	27 (7%)	15	47
All	All	866/928 (93%)	817 (94%)	49 (6%)	18	51

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	38	ARG
2	A	49	LEU
2	A	55	ARG
2	A	56	LEU
2	A	68	VAL
2	A	99	THR
2	A	115	ILE

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Mol	Chain	Res	Type
2	A	137	LYS
2	A	236	LEU
2	A	255	ARG
2	A	282	VAL
2	A	326	LEU
2	A	333	LEU
2	A	344	ARG
2	A	347	VAL
2	A	485	GLU
2	A	493	ILE
2	A	496	ASN
2	A	500	ILE
2	A	502	LEU
2	A	513	LEU
2	A	525	LEU
2	B	38	ARG
2	B	49	LEU
2	B	55	ARG
2	B	56	LEU
2	B	77	VAL
2	B	99	THR
2	B	102	LYS
2	B	180	THR
2	B	236	LEU
2	B	254	LEU
2	B	282	VAL
2	B	314	LEU
2	B	318	ASP
2	B	325	ASN
2	B	328	LEU
2	B	333	LEU
2	B	347	VAL
2	B	373	GLN
2	B	410	ARG
2	B	435	VAL
2	B	448	LYS
2	B	476	VAL
2	B	485	GLU
2	B	493	ILE
2	B	496	ASN
2	B	502	LEU
2	B	525	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	47	HIS
2	A	104	ASN
2	A	114	GLN
2	A	133	HIS
2	A	261	ASN
2	A	286	HIS
2	A	352	HIS
2	A	382	GLN
2	A	422	ASN
2	A	484	GLN
2	A	496	ASN
2	B	47	HIS
2	B	197	HIS
2	B	261	ASN
2	B	286	HIS
2	B	325	ASN
2	B	382	GLN
2	B	420	ASN
2	B	432	ASN
2	B	496	ASN
2	B	505	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	D	70/71 (98%)	15 (21%)	3 (4%)

All (15) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	D	906	G
1	D	908	U
1	D	909	G
1	D	910	G
1	D	917	G
1	D	918	G
1	D	919	C
1	D	932	U
1	D	933	U
1	D	935	A

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Mol	Chain	Res	Type
1	D	944	U
1	D	945	U
1	D	946	A
1	D	958	A
1	D	960	C

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	D	917	G
1	D	934	C
1	D	957	A

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	SEP	A	1002	-	9,10,10	1.59	1 (11%)	9,14,14	1.89	2 (22%)
3	SEP	A	1001	-	9,10,10	1.59	1 (11%)	9,14,14	1.79	2 (22%)

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	SEP	A	1002	-	-	1/10/10/10	-
3	SEP	A	1001	-	-	6/10/10/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1001	SEP	P-O1P	3.52	1.61	1.50
3	A	1002	SEP	P-O1P	3.51	1.61	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1002	SEP	OG-CB-CA	4.78	112.23	108.06
3	A	1001	SEP	OG-CB-CA	4.48	111.97	108.06
3	A	1002	SEP	O3P-P-OG	2.35	112.80	106.67
3	A	1001	SEP	O3P-P-OG	2.27	112.60	106.67

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1001	SEP	O-C-CA-CB
3	A	1001	SEP	CB-OG-P-O1P
3	A	1001	SEP	CB-OG-P-O2P
3	A	1001	SEP	CB-OG-P-O3P
3	A	1001	SEP	CA-CB-OG-P
3	A	1002	SEP	CA-CB-OG-P
3	A	1001	SEP	OXT-C-CA-N

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1002	SEP	1	0
3	A	1001	SEP	4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	D	71/71 (100%)	-0.01	0 100 100	104, 195, 200, 200	0
2	A	534/534 (100%)	-0.26	9 (1%) 69 49	20, 92, 197, 202	0
2	B	465/534 (87%)	-0.35	4 (0%) 81 64	25, 80, 185, 202	0
All	All	1070/1139 (93%)	-0.28	13 (1%) 76 58	20, 91, 198, 202	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	161	VAL	2.7
2	A	171	PHE	2.6
2	A	384	ILE	2.5
2	A	416	VAL	2.5
2	A	493	ILE	2.4
2	A	168	PHE	2.4
2	B	331	GLU	2.3
2	B	228	THR	2.3
2	A	83	PHE	2.2
2	B	388	ALA	2.2
2	B	493	ILE	2.2
2	A	380	ILE	2.1
2	A	529	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	SEP	A	1002	11/11	0.92	0.10	41,77,116,154	0
3	SEP	A	1001	11/11	0.95	0.09	41,103,137,172	0

6.5 Other polymers [i](#)

There are no such residues in this entry.