



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:51 PM GMT

PDB ID : 3V81
Title : Crystal structure of HIV-1 reverse transcriptase (RT) with DNA and the non-nucleoside inhibitor nevirapine
Authors : Das, K.; Martinez, S.E.; Arnold, E.
Deposited on : 2011-12-22
Resolution : 2.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

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

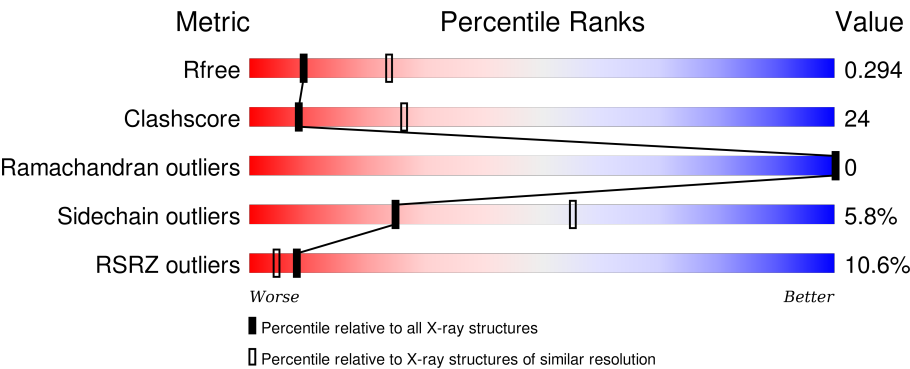
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	91344	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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	556	<div><div>14%</div><div>57%</div><div>40%</div><div>.</div></div>
1	C	556	<div><div>15%</div><div>54%</div><div>44%</div><div>.</div></div>
2	B	428	<div><div>4%</div><div>56%</div><div>36%</div><div>.</div><div>.</div></div>
2	D	428	<div><div>6%</div><div>60%</div><div>34%</div><div>.</div><div>.</div></div>
3	E	27	<div><div>7%</div><div>26%</div><div>33%</div><div>15%</div><div>26%</div></div>

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Mol	Chain	Length	Quality of chain
3	T	27	<div><div></div><div></div><div></div><div></div><div></div></div> <div>11%19%44%11%26%</div>
4	F	21	<div><div></div><div></div><div></div><div></div><div></div></div> <div>5%14%62%19%5%</div>
4	P	21	<div><div></div><div></div><div></div><div></div><div></div></div> <div>5%14%52%29%5%</div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17510 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 HIV-1 Reverse Transcriptase P66 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4512	2920	751	833	8			
1	C	555	Total	C	N	O	S	0	0	0
			4512	2920	751	833	8			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	EXPRESSION TAG	UNP P03366
A	0	VAL	-	EXPRESSION TAG	UNP P03366
A	258	CYS	GLN	ENGINEERED MUTATION	UNP P03366
A	280	SER	CYS	ENGINEERED MUTATION	UNP P03366
A	498	ASN	ASP	ENGINEERED MUTATION	UNP P03366
C	-1	MET	-	EXPRESSION TAG	UNP P03366
C	0	VAL	-	EXPRESSION TAG	UNP P03366
C	258	CYS	GLN	ENGINEERED MUTATION	UNP P03366
C	280	SER	CYS	ENGINEERED MUTATION	UNP P03366
C	498	ASN	ASP	ENGINEERED MUTATION	UNP P03366

- Molecule 2 is a protein called HIV-1 Reverse Transcriptase P51 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	412	Total	C	N	O	S	0	0	0
			3400	2212	563	619	6			
2	D	412	Total	C	N	O	S	0	0	0
			3400	2212	563	619	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	ENGINEERED MUTATION	UNP P03366
D	280	SER	CYS	ENGINEERED MUTATION	UNP P03366

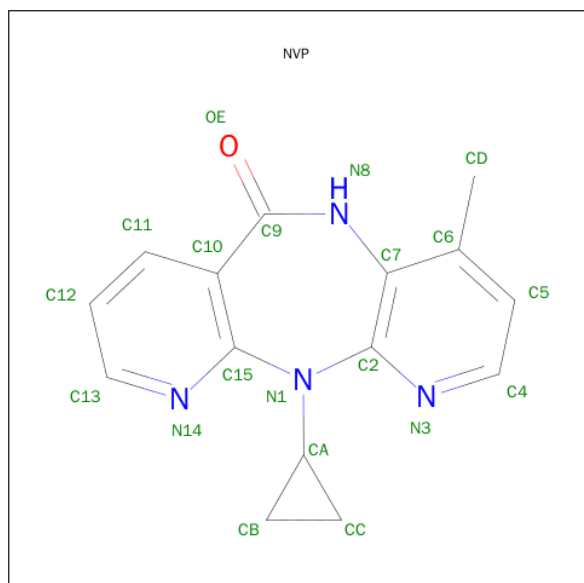
- Molecule 3 is a DNA chain called DNA (5'-D(*AP*TP*GP*GP*AP*AP*GP*GP*CP*GP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*GP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	20	Total	C	N	O	P	0	0	0
			415	194	85	116	20			
3	E	20	Total	C	N	O	P	0	0	0
			415	194	85	116	20			

- Molecule 4 is a DNA chain called DNA (5'-D(*A*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MRG)P*CP*GP*CP*CP*(ATM))-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	P	20	Total	C	N	O	P	S	0	0	0
			408	195	72	121	19	1			
4	F	20	Total	C	N	O	P	S	0	0	0
			408	195	72	121	19	1			

- Molecule 5 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).



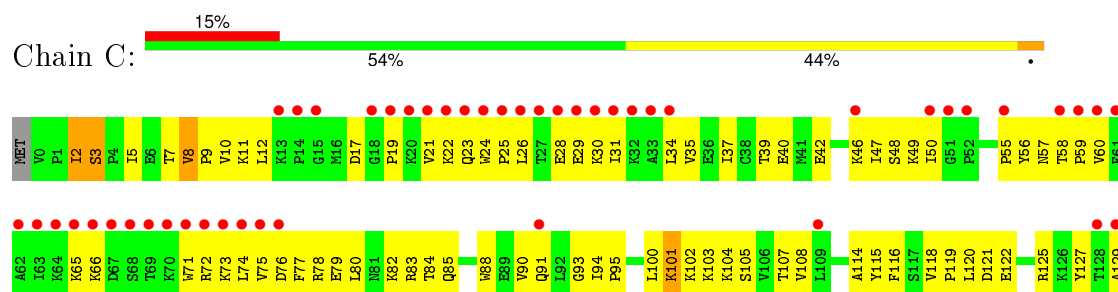
3 Residue-property plots

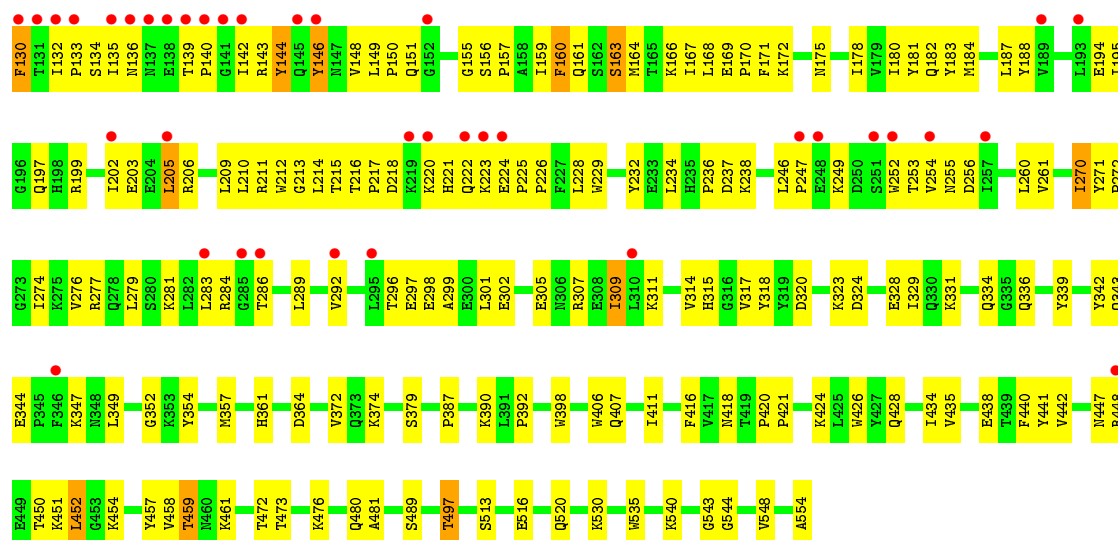
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: HIV-1 Reverse Transcriptase P66 subunit

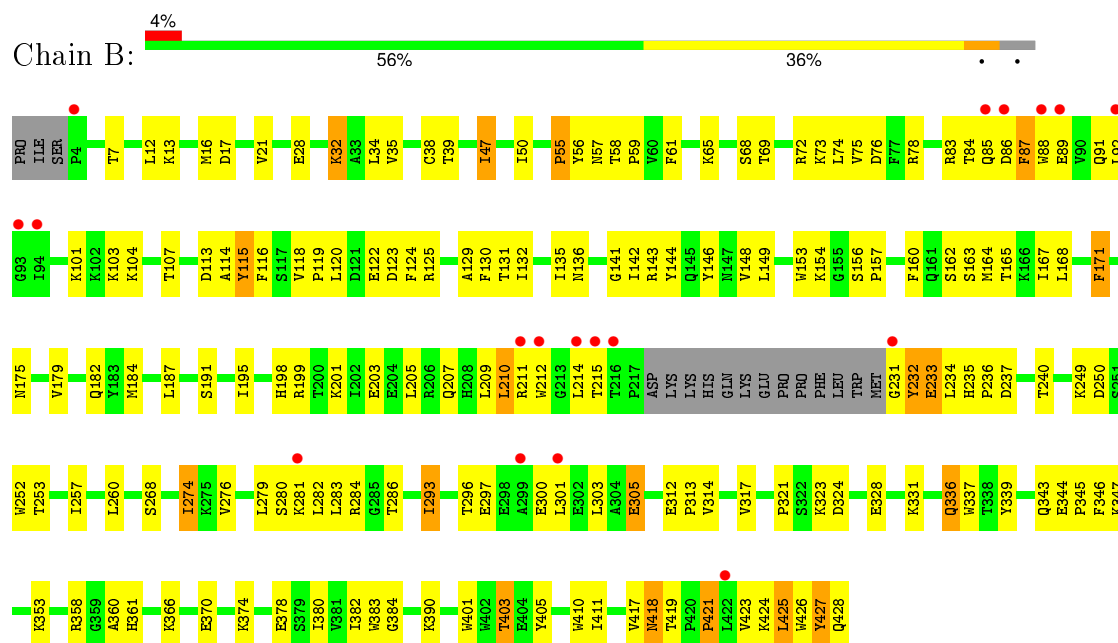


• Molecule 1: HIV-1 Reverse Transcriptase P66 subunit

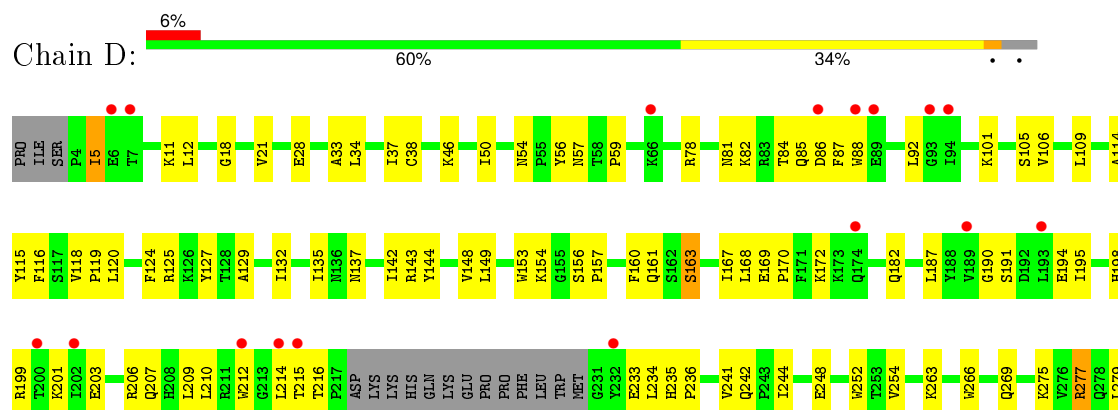


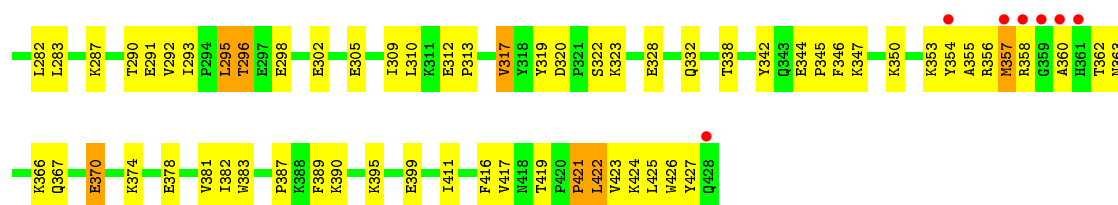


● Molecule 2: HIV-1 Reverse Transcriptase P51 subunit

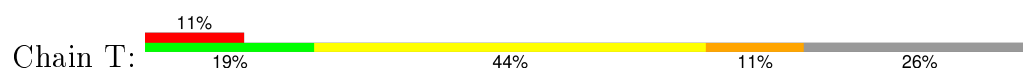


● Molecule 2: HIV-1 Reverse Transcriptase P51 subunit

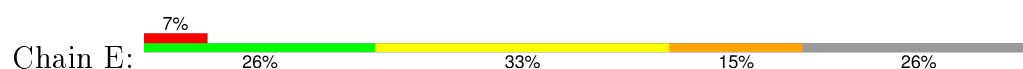




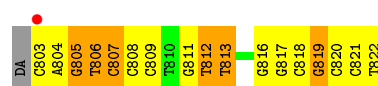
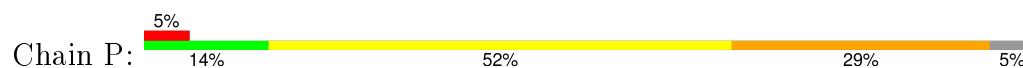
- Molecule 3: DNA (5'-D(*AP*TP*GP*GP*AP*AP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*GP*TP*G)-3')



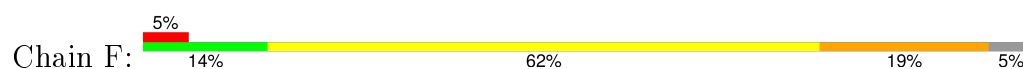
- Molecule 3: DNA (5'-D(*AP*TP*GP*GP*AP*AP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*GP*TP*G)-3')



- Molecule 4: DNA (5'-D(*A*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MR G)P*CP*GP*CP*CP*(ATM))-3')



- Molecule 4: DNA (5'-D(*A*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*(MR G)P*CP*GP*CP*CP*(ATM))-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.82Å 132.05Å 142.73Å 90.00° 100.84° 90.00°	Depositor
Resolution (Å)	38.41 – 2.85 38.41 – 2.85	Depositor EDS
% Data completeness (in resolution range)	98.7 (38.41-2.85) 98.6 (38.41-2.85)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.251 , 0.298 0.248 , 0.294	Depositor DCC
R_{free} test set	2288 reflections (3.13%)	DCC
Wilson B-factor (Å ²)	54.2	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 62.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 75371 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	17510	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ATM, MRG, 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	0.49	0/4630	0.61	0/6290
1	C	0.49	0/4630	0.63	0/6290
2	B	0.51	0/3497	0.66	2/4751 (0.0%)
2	D	0.48	0/3497	0.63	0/4751
3	E	0.70	0/467	1.55	9/719 (1.3%)
3	T	0.67	0/467	1.46	6/719 (0.8%)
4	F	0.72	0/400	1.66	13/612 (2.1%)
4	P	1.15	2/400 (0.5%)	1.74	13/612 (2.1%)
All	All	0.53	2/17988 (0.0%)	0.79	43/24744 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1
2	D	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	805	DG	O3'-P	-14.74	1.43	1.61
4	P	806	DT	O3'-P	-9.82	1.49	1.61

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	807	DC	O4'-C1'-N1	12.19	116.53	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	807	DC	O4'-C1'-N1	10.73	115.51	108.00
3	T	719	DG	O4'-C1'-N9	10.60	115.42	108.00
4	P	820	DC	O4'-C4'-C3'	-8.80	100.72	106.00
4	P	805	DG	O3'-P-O5'	8.18	119.54	104.00
4	F	820	DC	O4'-C4'-C3'	-8.11	101.13	106.00
3	E	719	DG	O4'-C1'-N9	7.92	113.55	108.00
3	T	717	DC	O4'-C1'-N1	7.55	113.29	108.00
3	E	709	DC	O4'-C4'-C3'	-7.55	101.47	106.00
3	E	722	DA	O4'-C1'-N9	7.34	113.14	108.00
3	T	709	DC	O4'-C4'-C3'	-7.27	101.59	104.50
4	F	821	DC	O4'-C4'-C3'	-7.00	101.70	104.50
4	P	809	DC	O4'-C1'-N1	6.90	112.83	108.00
3	E	719	DG	C1'-O4'-C4'	-6.81	103.29	110.10
4	P	818	DC	P-O3'-C3'	6.73	127.78	119.70
4	F	814	DC	O4'-C1'-N1	6.45	112.51	108.00
4	F	813	DT	C5-C4-O4	-6.27	120.51	124.90
4	F	816	DG	O4'-C1'-N9	6.13	112.29	108.00
3	E	719	DG	O4'-C1'-C2'	-6.08	101.03	105.90
3	E	716	DA	O4'-C1'-N9	5.93	112.15	108.00
4	P	813	DT	C5-C4-O4	-5.91	120.76	124.90
4	F	813	DT	O4'-C1'-C2'	-5.88	101.19	105.90
4	F	812	DT	N3-C4-O4	5.57	123.24	119.90
3	T	725	DG	C1'-O4'-C4'	-5.56	104.54	110.10
4	P	812	DT	O4'-C1'-N1	-5.47	104.17	108.00
2	B	425	LEU	CA-CB-CG	-5.45	102.77	115.30
3	E	715	DA	O4'-C1'-C2'	-5.44	101.55	105.90
3	E	723	DC	C1'-O4'-C4'	-5.41	104.69	110.10
3	T	721	DG	C4'-C3'-C2'	-5.37	98.27	103.10
4	F	813	DT	O4'-C1'-N1	5.33	111.73	108.00
4	F	810	DT	O4'-C1'-C2'	-5.30	101.66	105.90
3	T	725	DG	O4'-C1'-N9	5.25	111.67	108.00
4	P	805	DG	OP2-P-O3'	-5.24	93.67	105.20
4	P	805	DG	P-O3'-C3'	5.17	125.90	119.70
4	F	811	DG	O4'-C1'-N9	5.15	111.60	108.00
2	B	401	TRP	N-CA-C	5.14	124.88	111.00
4	P	820	DC	O4'-C1'-N1	5.08	111.56	108.00
4	F	813	DT	N3-C4-O4	5.07	122.94	119.90
4	F	814	DC	O4'-C1'-C2'	-5.06	101.85	105.90
3	E	708	DG	O4'-C1'-N9	5.04	111.53	108.00
4	P	809	DC	O4'-C4'-C3'	5.04	109.02	106.00
4	P	819	DG	O4'-C4'-C3'	-5.03	102.49	104.50
4	P	818	DC	OP1-P-O3'	5.03	116.26	105.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	421	PRO	Peptide
2	D	421	PRO	Peptide

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	4512	0	4570	228	0
1	C	4512	0	4570	256	0
2	B	3400	0	3433	158	0
2	D	3400	0	3433	143	0
3	E	415	0	222	12	0
3	T	415	0	222	21	0
4	F	408	0	231	18	0
4	P	408	0	231	23	0
5	A	20	0	14	2	0
5	C	20	0	14	1	0
All	All	17510	0	16940	827	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 24.

All (827) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:ILE:HG22	1:C:212:TRP:HD1	1.09	1.17
1:A:5:ILE:HG22	1:A:212:TRP:HD1	1.17	1.08
1:C:21:VAL:H	1:C:57:ASN:HB2	0.93	1.06
1:A:283:LEU:HD13	4:P:817:MRG:H231	1.39	1.03
1:C:283:LEU:HD23	4:F:817:MRG:H231	1.44	0.99
1:A:195:ILE:HG22	1:A:199:ARG:HH12	1.24	0.99
1:C:21:VAL:N	1:C:57:ASN:HB2	1.77	0.98
1:C:5:ILE:HG22	1:C:212:TRP:CD1	2.00	0.96
2:B:115:TYR:HD1	2:B:156:SER:HB3	1.30	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:282:LEU:HD21	2:D:296:THR:HG22	1.48	0.95
1:C:155:GLY:O	1:C:159:ILE:HG13	1.66	0.95
1:A:406:TRP:CZ3	1:A:407:GLN:HG3	2.01	0.94
1:A:261:VAL:HG13	1:A:276:VAL:HG21	1.47	0.94
1:A:5:ILE:HG22	1:A:212:TRP:CD1	2.04	0.92
2:D:88:TRP:CE2	2:D:154:LYS:HG2	2.06	0.90
2:B:107:THR:HG23	2:B:232:TYR:HB2	1.55	0.89
1:A:120:LEU:HD12	1:A:121:ASP:H	1.38	0.89
2:B:28:GLU:HB2	2:B:135:ILE:HD11	1.54	0.88
2:D:168:LEU:HB3	2:D:172:LYS:HE3	1.52	0.88
1:C:261:VAL:HG13	1:C:276:VAL:HG21	1.56	0.88
2:B:282:LEU:HD21	2:B:296:THR:HG23	1.57	0.87
1:C:447:ASN:ND2	1:C:450:THR:H	1.73	0.86
1:A:2:ILE:HD12	1:A:2:ILE:N	1.90	0.86
2:D:421:PRO:HB2	2:D:423:VAL:H	1.41	0.85
1:A:114:ALA:HB1	1:A:160:PHE:HE1	1.42	0.84
1:C:125:ARG:HG2	1:C:146:TYR:O	1.77	0.84
1:C:2:ILE:HD12	1:C:2:ILE:N	1.91	0.84
4:F:816:DG:H2''	4:F:817:MRG:H5'	1.59	0.84
1:A:135:ILE:HG23	1:A:136:ASN:H	1.40	0.84
1:C:21:VAL:H	1:C:57:ASN:CB	1.86	0.84
1:C:2:ILE:HD12	1:C:2:ILE:H	1.41	0.84
1:C:438:GLU:OE2	1:C:459:THR:HG21	1.76	0.83
1:A:32:LYS:HD3	1:A:135:ILE:HB	1.60	0.83
1:A:155:GLY:O	1:A:159:ILE:HG13	1.78	0.83
1:C:296:THR:HB	1:C:299:ALA:HB3	1.59	0.83
1:A:24:TRP:CD2	1:A:25:PRO:HD2	2.13	0.83
2:B:164:MET:HE2	2:B:168:LEU:HD11	1.58	0.83
2:D:163:SER:O	2:D:167:ILE:HG22	1.79	0.82
2:D:115:TYR:CD1	2:D:156:SER:HB3	2.14	0.82
1:A:34:LEU:HD22	1:A:73:LYS:HE2	1.59	0.81
1:C:139:THR:HB	1:C:140:PRO:HD2	1.61	0.81
2:B:115:TYR:CD1	2:B:156:SER:HB3	2.15	0.81
1:A:195:ILE:HG22	1:A:199:ARG:NH1	1.94	0.80
1:A:281:LYS:HE3	1:A:284:ARG:HH12	1.47	0.80
1:C:139:THR:HB	1:C:140:PRO:CD	2.12	0.79
4:P:804:DA:H2''	4:P:805:DG:H5'	1.64	0.79
1:C:398:TRP:CZ2	1:C:411:ILE:HG13	2.17	0.79
1:C:77:PHE:CD2	1:C:80:LEU:HB3	2.17	0.79
1:C:447:ASN:HD22	1:C:450:THR:H	1.29	0.79
4:P:816:DG:H2''	4:P:817:MRG:H5'	1.63	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:ILE:HD13	1:A:530:LYS:HB3	1.66	0.78
1:A:41:MET:HB3	1:A:46:LYS:HB2	1.65	0.78
1:C:21:VAL:HG12	1:C:22:LYS:H	1.48	0.77
2:B:142:ILE:HG22	2:B:144:TYR:HE2	1.49	0.77
2:B:167:ILE:HG23	2:B:212:TRP:CD1	2.20	0.77
2:B:91:GLN:HG2	2:B:92:LEU:HG	1.67	0.77
1:C:223:LYS:O	1:C:225:PRO:HD3	1.85	0.77
1:C:435:VAL:HG22	2:D:290:THR:HG21	1.66	0.76
2:B:314:VAL:O	2:B:317:VAL:HG22	1.84	0.76
1:A:24:TRP:CG	1:A:25:PRO:HD2	2.20	0.76
2:B:73:LYS:HE2	2:B:146:TYR:OH	1.86	0.76
1:C:372:VAL:HG11	1:C:411:ILE:HG23	1.67	0.75
2:B:142:ILE:HG22	2:B:144:TYR:CE2	2.21	0.75
3:T:721:DG:H2''	3:T:722:DA:C8	2.22	0.75
2:D:254:VAL:HG13	2:D:283:LEU:HD22	1.67	0.75
1:A:221:HIS:CG	1:A:224:GLU:HB2	2.22	0.74
1:C:31:ILE:HD12	1:C:135:ILE:HG13	1.67	0.74
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.22	0.74
2:D:115:TYR:HD1	2:D:156:SER:HB3	1.52	0.74
2:B:28:GLU:CB	2:B:135:ILE:HD11	2.16	0.74
1:C:199:ARG:HH21	1:C:223:LYS:HD3	1.51	0.74
1:A:221:HIS:HB2	1:A:224:GLU:HB2	1.69	0.74
1:A:406:TRP:CH2	1:A:407:GLN:HG3	2.23	0.73
2:D:266:TRP:O	2:D:269:GLN:HG3	1.88	0.73
2:B:366:LYS:HG3	2:B:405:TYR:CD1	2.23	0.73
1:A:271:TYR:CE1	1:A:314:VAL:HG12	2.23	0.73
4:P:804:DA:H2'	4:P:805:DG:C8	2.24	0.73
1:C:195:ILE:HD11	1:C:223:LYS:HD3	1.71	0.73
1:C:206:ARG:HG2	1:C:216:THR:OG1	1.88	0.72
2:D:277:ARG:H	2:D:277:ARG:HD2	1.54	0.72
1:A:125:ARG:HG2	1:A:146:TYR:O	1.89	0.72
1:C:274:ILE:HD12	1:C:309:ILE:HD11	1.72	0.72
1:C:49:LYS:O	1:C:50:ILE:HD12	1.89	0.72
1:A:406:TRP:CH2	2:B:418:ASN:HA	2.25	0.72
2:B:87:PHE:O	2:B:91:GLN:HB3	1.88	0.72
2:D:86:ASP:OD1	2:D:87:PHE:N	2.23	0.72
1:A:60:VAL:HG21	1:A:132:ILE:HD12	1.71	0.71
2:B:124:PHE:CE2	2:B:153:TRP:CZ2	2.78	0.71
3:E:720:DG:H4'	3:E:721:DG:OP1	1.91	0.71
2:B:171:PHE:HE2	2:B:205:LEU:HB2	1.56	0.71
1:A:167:ILE:O	1:A:170:PRO:HD2	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:HIS:HD1	1:A:513:SER:HG	1.34	0.71
1:A:500:GLN:NE2	3:T:722:DA:OP1	2.24	0.70
2:B:282:LEU:HB3	2:B:293:ILE:HD11	1.73	0.70
2:D:120:LEU:HD23	2:D:125:ARG:HG2	1.72	0.70
2:B:182:GLN:HG3	2:B:187:LEU:HD12	1.73	0.70
4:P:807:DC:H4'	4:P:808:DC:OP1	1.92	0.70
2:B:13:LYS:HG3	2:B:16:MET:HE1	1.73	0.70
1:C:17:ASP:O	1:C:83:ARG:HD3	1.92	0.70
1:C:77:PHE:HD2	1:C:80:LEU:HB3	1.58	0.69
2:B:56:TYR:O	2:B:143:ARG:NH2	2.24	0.69
2:B:323:LYS:HB2	2:B:343:GLN:HE21	1.56	0.69
2:D:421:PRO:HB2	2:D:423:VAL:N	2.07	0.69
3:E:719:DG:H4'	3:E:720:DG:OP1	1.92	0.69
1:C:296:THR:HB	1:C:299:ALA:CB	2.22	0.69
1:A:288:ALA:HB3	1:A:291:GLU:HG3	1.74	0.69
1:C:75:VAL:HG12	1:C:77:PHE:HD1	1.57	0.69
1:A:2:ILE:HD12	1:A:2:ILE:H	1.57	0.68
2:B:21:VAL:HG13	2:B:59:PRO:HD3	1.75	0.68
2:D:263:LYS:HA	2:D:425:LEU:HD23	1.75	0.68
2:B:154:LYS:HG2	2:B:184:MET:SD	2.33	0.68
1:C:171:PHE:CE1	1:C:205:LEU:HB2	2.28	0.68
2:D:92:LEU:HD21	2:D:161:GLN:OE1	1.94	0.68
4:F:807:DC:H2''	4:F:808:DC:O5'	1.93	0.68
2:D:423:VAL:HG13	2:D:426:TRP:CD1	2.28	0.68
1:A:172:LYS:HE2	1:A:180:ILE:HD12	1.74	0.68
2:D:182:GLN:HB2	2:D:187:LEU:HD12	1.76	0.68
1:A:77:PHE:O	1:A:80:LEU:N	2.27	0.68
1:C:459:THR:HG22	1:C:461:LYS:N	2.10	0.67
1:A:26:LEU:HD22	1:A:60:VAL:HG11	1.76	0.67
3:E:717:DC:H2'	3:E:718:DA:C8	2.29	0.67
2:B:122:GLU:HA	2:B:125:ARG:HD2	1.76	0.67
1:C:281:LYS:HG2	1:C:284:ARG:NH1	2.09	0.67
1:C:60:VAL:HG23	1:C:130:PHE:HE1	1.60	0.67
2:B:323:LYS:HB2	2:B:343:GLN:NE2	2.09	0.67
1:A:103:LYS:O	1:A:236:PRO:HB3	1.93	0.67
1:A:156:SER:HB2	1:A:157:PRO:HD3	1.76	0.66
2:B:214:LEU:HD12	2:B:215:THR:H	1.60	0.66
2:D:358:ARG:HG2	2:D:366:LYS:HD3	1.75	0.66
1:A:238:LYS:HD2	1:A:315:HIS:ND1	2.10	0.66
2:B:249:LYS:HB2	2:B:252:TRP:CE2	2.30	0.66
3:E:714:DG:H2''	3:E:715:DA:C8	2.31	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:281:LYS:HG2	1:C:284:ARG:HH12	1.60	0.66
2:D:88:TRP:CD2	2:D:154:LYS:HG2	2.31	0.66
3:T:706:DA:N1	4:P:822:ATM:O4	2.28	0.66
1:C:47:ILE:HG22	1:C:48:SER:N	2.11	0.66
1:C:29:GLU:OE2	1:C:71:TRP:HH2	1.77	0.66
1:C:271:TYR:CE1	1:C:314:VAL:HG12	2.31	0.66
2:D:312:GLU:HB3	2:D:313:PRO:HD2	1.76	0.65
1:A:542:ILE:HG23	2:B:283:LEU:HD13	1.76	0.65
1:A:21:VAL:O	1:A:57:ASN:ND2	2.29	0.65
1:A:171:PHE:CD1	1:A:205:LEU:HD12	2.30	0.65
1:A:5:ILE:CG2	1:A:212:TRP:HD1	2.03	0.65
2:B:47:ILE:HD12	2:B:144:TYR:CD1	2.32	0.65
2:D:320:ASP:OD1	2:D:322:SER:HB3	1.96	0.65
1:A:305:GLU:O	1:A:309:ILE:HG13	1.95	0.64
2:B:73:LYS:NZ	2:B:130:PHE:CZ	2.65	0.64
2:B:317:VAL:HG12	2:B:347:LYS:HB3	1.79	0.64
1:C:441:TYR:CD1	1:C:544:GLY:HA3	2.33	0.64
1:C:120:LEU:HG	1:C:121:ASP:H	1.61	0.64
2:B:374:LYS:O	2:B:378:GLU:HG3	1.97	0.64
1:A:115:TYR:HD2	1:A:151:GLN:HE22	1.46	0.64
2:B:114:ALA:HB2	2:B:214:LEU:HD22	1.78	0.64
1:C:148:VAL:O	1:C:150:PRO:HD3	1.98	0.64
1:C:457:TYR:HA	1:C:548:VAL:HG21	1.78	0.64
2:B:425:LEU:HD22	2:B:425:LEU:N	2.12	0.64
2:B:104:LYS:HA	2:B:237:ASP:OD1	1.97	0.64
1:C:459:THR:HG22	1:C:461:LYS:H	1.61	0.64
1:C:2:ILE:HG22	1:C:3:SER:H	1.62	0.64
2:B:28:GLU:HB2	2:B:135:ILE:CD1	2.28	0.63
1:A:164:MET:O	1:A:168:LEU:HD13	1.97	0.63
1:A:221:HIS:CB	1:A:224:GLU:HB2	2.27	0.63
1:C:473:THR:OG1	1:C:476:LYS:HG3	1.98	0.63
2:B:370:GLU:O	2:B:374:LYS:HG3	1.98	0.63
2:D:395:LYS:HE3	2:D:399:GLU:OE2	1.99	0.63
1:A:115:TYR:CD1	1:A:160:PHE:HD1	2.17	0.63
2:B:74:LEU:HD12	2:B:75:VAL:N	2.14	0.63
2:B:171:PHE:HD2	2:B:205:LEU:HD13	1.64	0.63
1:A:46:LYS:HE3	1:A:116:PHE:HB3	1.81	0.62
1:C:450:THR:OG1	1:C:452:LEU:HB2	1.99	0.62
1:A:225:PRO:HA	1:A:226:PRO:C	2.20	0.62
1:A:88:TRP:CZ2	2:B:57:ASN:HB2	2.33	0.62
1:C:30:LYS:HD2	1:C:30:LYS:N	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:214:LEU:HD12	2:D:215:THR:H	1.64	0.62
1:C:199:ARG:NH2	1:C:223:LYS:HD3	2.14	0.62
1:C:156:SER:HB2	1:C:157:PRO:HD3	1.81	0.62
2:D:358:ARG:HB3	2:D:370:GLU:OE1	1.98	0.62
1:C:175:ASN:HB3	1:C:178:ILE:HD13	1.81	0.62
1:C:296:THR:CG2	1:C:298:GLU:HG2	2.29	0.62
2:B:124:PHE:HE2	2:B:153:TRP:CZ2	2.18	0.62
1:A:56:TYR:HB2	1:A:129:ALA:HB3	1.81	0.61
1:A:88:TRP:HB2	2:B:55:PRO:HA	1.80	0.61
4:F:822:ATM:H6	4:F:822:ATM:OP1	1.99	0.61
1:A:406:TRP:HH2	2:B:418:ASN:HA	1.65	0.61
2:D:101:LYS:O	2:D:236:PRO:HB2	1.99	0.61
2:B:101:LYS:O	2:B:236:PRO:HB2	2.00	0.61
2:B:382:ILE:HG22	2:B:383:TRP:CD2	2.35	0.61
2:B:209:LEU:HD23	2:B:214:LEU:HD23	1.81	0.61
2:D:421:PRO:HB3	2:D:423:VAL:HB	1.82	0.61
1:A:404:GLU:OE1	1:A:404:GLU:HA	2.01	0.61
1:C:459:THR:CG2	1:C:461:LYS:H	2.14	0.61
2:B:142:ILE:CG2	2:B:144:TYR:HE2	2.13	0.61
1:A:142:ILE:HG13	1:A:143:ARG:H	1.65	0.61
1:A:175:ASN:HB3	1:A:178:ILE:HG12	1.82	0.61
1:C:160:PHE:C	1:C:160:PHE:CD2	2.73	0.61
3:T:720:DG:H2''	3:T:721:DG:C5'	2.30	0.61
1:C:344:GLU:HB2	1:C:347:LYS:HB2	1.83	0.61
1:A:50:ILE:HG22	1:A:52:PRO:HD2	1.81	0.61
2:B:297:GLU:O	2:B:301:LEU:HG	2.01	0.61
1:A:224:GLU:N	1:A:225:PRO:HD2	2.16	0.61
1:A:172:LYS:CE	1:A:180:ILE:HD12	2.31	0.61
2:D:354:TYR:CE2	2:D:374:LYS:HB3	2.36	0.61
1:A:483:TYR:CZ	1:A:487:GLN:NE2	2.69	0.61
1:A:434:ILE:CD1	1:A:530:LYS:HB3	2.31	0.60
3:T:716:DA:H2'	3:T:717:DC:H6	1.65	0.60
2:B:268:SER:HB3	2:B:274:ILE:HB	1.83	0.60
2:B:419:THR:HG22	2:B:419:THR:O	2.00	0.60
1:C:58:THR:HB	1:C:77:PHE:HB3	1.83	0.60
1:C:167:ILE:O	1:C:170:PRO:HD2	2.01	0.60
3:T:720:DG:H4'	3:T:721:DG:OP1	2.02	0.60
1:C:210:LEU:O	1:C:213:GLY:N	2.27	0.60
1:C:90:VAL:HB	1:C:161:GLN:OE1	2.01	0.60
1:A:543:GLY:HA2	2:B:283:LEU:O	2.01	0.60
1:C:441:TYR:CG	1:C:544:GLY:HA3	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:LEU:CG	1:C:121:ASP:H	2.14	0.60
1:C:379:SER:CB	1:C:387:PRO:HD3	2.31	0.60
1:A:120:LEU:HD12	1:A:121:ASP:N	2.13	0.60
2:D:360:ALA:HA	2:D:367:GLN:NE2	2.17	0.60
1:C:281:LYS:O	1:C:284:ARG:HG3	2.02	0.60
2:B:84:THR:O	2:B:87:PHE:HB3	2.01	0.60
1:A:301:LEU:O	1:A:305:GLU:HG3	2.02	0.60
1:C:434:ILE:HD13	1:C:530:LYS:HB3	1.83	0.60
2:D:33:ALA:O	2:D:37:ILE:HG13	2.02	0.59
3:E:720:DG:H2''	3:E:721:DG:O5'	2.02	0.59
3:T:716:DA:H2''	3:T:717:DC:O5'	2.02	0.59
2:D:421:PRO:CB	2:D:423:VAL:HB	2.32	0.59
2:D:423:VAL:HG13	2:D:426:TRP:NE1	2.17	0.59
1:C:220:LYS:HE2	1:C:222:GLN:HG2	1.84	0.59
1:A:548:VAL:HA	1:A:551:LEU:HD12	1.84	0.59
3:T:706:DA:N1	4:P:822:ATM:C4	2.65	0.59
1:C:160:PHE:HD2	1:C:160:PHE:C	2.06	0.59
1:A:2:ILE:CD1	1:A:2:ILE:N	2.61	0.59
1:A:26:LEU:CD2	1:A:132:ILE:HD11	2.33	0.59
1:A:171:PHE:CE1	1:A:205:LEU:HD12	2.37	0.59
1:C:206:ARG:HH22	1:C:218:ASP:HB2	1.68	0.59
1:A:50:ILE:HG21	1:A:53:GLU:HG2	1.83	0.59
2:D:28:GLU:CB	2:D:135:ILE:HD11	2.32	0.59
1:C:47:ILE:HG21	1:C:144:TYR:HD1	1.68	0.59
2:B:390:LYS:HB3	2:B:417:VAL:HG21	1.84	0.59
1:C:473:THR:HG21	4:F:809:DC:P	2.43	0.59
1:A:17:ASP:OD1	1:A:18:GLY:N	2.36	0.58
1:A:128:THR:HG21	1:A:146:TYR:HB2	1.85	0.58
2:B:171:PHE:CE2	2:B:205:LEU:HB2	2.38	0.58
1:A:259:LYS:HG3	4:P:819:DG:OP1	2.03	0.58
1:C:163:SER:HA	1:C:166:LYS:HE3	1.84	0.58
1:A:58:THR:HG23	1:A:129:ALA:O	2.04	0.58
2:D:28:GLU:HB2	2:D:135:ILE:HD11	1.84	0.58
1:C:73:LYS:HG2	1:C:74:LEU:N	2.18	0.58
1:C:296:THR:HG21	1:C:298:GLU:HG2	1.85	0.58
1:A:77:PHE:HD2	1:A:80:LEU:HD23	1.67	0.58
2:B:297:GLU:HA	2:B:300:GLU:HB3	1.85	0.58
2:D:374:LYS:O	2:D:378:GLU:HG3	2.02	0.58
2:D:115:TYR:HE1	2:D:157:PRO:N	2.02	0.58
1:C:543:GLY:HA3	2:D:283:LEU:O	2.03	0.58
1:C:218:ASP:OD2	1:C:222:GLN:NE2	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:118:VAL:HG12	2:D:119:PRO:O	2.03	0.58
1:C:361:HIS:ND1	1:C:513:SER:OG	2.25	0.58
1:C:108:VAL:HG22	1:C:188:TYR:CD1	2.39	0.58
1:A:159:ILE:O	1:A:163:SER:OG	2.22	0.58
1:C:59:PRO:HG2	1:C:76:ASP:HB3	1.85	0.58
3:T:721:DG:H2"	3:T:722:DA:H8	1.69	0.58
4:P:807:DC:H2"	4:P:808:DC:C5'	2.34	0.58
2:D:88:TRP:CD1	2:D:154:LYS:HB3	2.39	0.57
1:A:132:ILE:HG12	1:A:133:PRO:HD2	1.86	0.57
2:D:290:THR:HG22	2:D:290:THR:O	2.04	0.57
1:C:7:THR:HG21	1:C:122:GLU:OE2	2.04	0.57
1:C:361:HIS:HD1	1:C:513:SER:HG	1.48	0.57
1:A:194:GLU:HG3	1:A:197:GLN:HB3	1.87	0.57
1:C:277:ARG:O	1:C:277:ARG:HD3	2.04	0.57
1:A:281:LYS:HE3	1:A:284:ARG:NH1	2.18	0.57
1:C:164:MET:HE2	1:C:168:LEU:HD11	1.85	0.57
1:A:295:LEU:HB3	1:A:300:GLU:HG3	1.87	0.57
2:D:114:ALA:N	2:D:214:LEU:HD13	2.19	0.57
4:P:806:DT:H2"	4:P:807:DC:O5'	2.05	0.57
1:C:328:GLU:HG3	1:C:390:LYS:HB2	1.87	0.57
3:T:706:DA:O5'	3:T:706:DA:H8	1.88	0.57
2:D:338:THR:HG21	2:D:427:TYR:HB2	1.86	0.57
1:A:34:LEU:CD2	1:A:73:LYS:HE2	2.33	0.56
1:C:135:ILE:HG12	1:C:136:ASN:OD1	2.05	0.56
1:A:2:ILE:HG22	1:A:3:SER:H	1.70	0.56
1:A:115:TYR:HD2	1:A:151:GLN:NE2	2.02	0.56
1:A:47:ILE:HG22	1:A:48:SER:N	2.20	0.56
1:A:372:VAL:HG11	1:A:411:ILE:HG23	1.87	0.56
2:B:231:GLY:C	2:B:233:GLU:H	2.08	0.56
2:D:382:ILE:HG22	2:D:383:TRP:CD2	2.40	0.56
1:C:324:ASP:O	1:C:343:GLN:HG2	2.05	0.56
3:T:720:DG:H2"	3:T:721:DG:H5"	1.85	0.56
1:A:77:PHE:CD2	1:A:80:LEU:HD23	2.40	0.56
2:B:253:THR:O	2:B:257:ILE:HG12	2.05	0.56
1:A:77:PHE:HB3	1:A:80:LEU:HB3	1.87	0.56
2:D:282:LEU:CD2	2:D:296:THR:HG22	2.31	0.56
1:C:184:MET:CE	4:F:822:ATM:H2'	2.35	0.56
1:A:276:VAL:HG22	1:A:276:VAL:O	2.06	0.56
1:A:160:PHE:HE2	1:A:164:MET:CE	2.19	0.56
1:C:435:VAL:HG22	2:D:290:THR:CG2	2.35	0.56
1:A:361:HIS:ND1	1:A:513:SER:OG	2.33	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:447:ASN:ND2	1:C:450:THR:HG23	2.21	0.56
1:C:130:PHE:CD2	1:C:144:TYR:CE2	2.94	0.56
1:A:253:THR:HG22	1:A:256:ASP:OD2	2.06	0.56
2:B:17:ASP:O	2:B:83:ARG:HD3	2.05	0.56
2:D:115:TYR:HE1	2:D:156:SER:C	2.08	0.55
2:D:332:GLN:HG3	2:D:338:THR:HG23	1.87	0.55
1:C:206:ARG:HH21	1:C:217:PRO:C	2.10	0.55
1:A:12:LEU:HD22	1:A:83:ARG:O	2.05	0.55
1:C:457:TYR:C	1:C:457:TYR:CD1	2.80	0.55
2:D:142:ILE:HG22	2:D:144:TYR:CE2	2.41	0.55
1:A:365:VAL:HG11	1:A:401:TRP:CD1	2.42	0.55
1:A:19:PRO:O	1:A:56:TYR:HA	2.06	0.55
2:D:203:GLU:HA	2:D:206:ARG:HB2	1.89	0.55
1:A:455:ALA:HB2	1:A:469:LEU:HD11	1.87	0.55
1:C:118:VAL:HG12	1:C:149:LEU:HG	1.88	0.55
2:B:88:TRP:CZ2	2:B:154:LYS:HD2	2.41	0.55
1:A:221:HIS:CD2	1:A:224:GLU:HB2	2.42	0.55
1:C:272:PRO:O	1:C:309:ILE:HD12	2.07	0.55
1:C:329:ILE:O	1:C:392:PRO:HD3	2.07	0.55
2:D:422:LEU:HD23	2:D:422:LEU:C	2.28	0.55
2:D:124:PHE:CE2	2:D:153:TRP:CZ2	2.94	0.55
1:A:160:PHE:HD2	1:A:160:PHE:C	2.11	0.55
4:P:807:DC:H2'	4:P:808:DC:C6	2.42	0.55
2:B:336:GLN:NE2	2:B:427:TYR:OH	2.40	0.55
2:B:162:SER:O	2:B:165:THR:HG22	2.06	0.55
1:A:226:PRO:HD3	1:A:235:HIS:CD2	2.42	0.54
1:A:90:VAL:HG11	1:A:161:GLN:HB3	1.89	0.54
1:C:47:ILE:CG2	1:C:48:SER:N	2.71	0.54
3:E:714:DG:H2''	3:E:715:DA:H8	1.72	0.54
3:T:711:DC:H2'	3:T:712:DC:C6	2.43	0.54
2:D:320:ASP:OD2	2:D:323:LYS:HE2	2.07	0.54
2:D:305:GLU:O	2:D:309:ILE:HG13	2.07	0.54
2:B:38:CYS:SG	2:B:132:ILE:HD11	2.47	0.54
1:C:120:LEU:HG	1:C:121:ASP:N	2.22	0.54
1:A:5:ILE:O	1:A:119:PRO:HG3	2.08	0.54
2:D:114:ALA:HB1	2:D:160:PHE:CZ	2.43	0.54
2:B:114:ALA:N	2:B:214:LEU:HD13	2.23	0.54
1:C:458:VAL:HG23	1:C:548:VAL:HB	1.89	0.54
1:A:118:VAL:HB	1:A:149:LEU:HD12	1.90	0.54
1:C:23:GLN:HE22	1:C:133:PRO:HD2	1.72	0.54
1:C:77:PHE:CE2	1:C:80:LEU:HD23	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:720:DG:H2''	3:T:721:DG:O5'	2.07	0.54
1:C:228:LEU:HA	1:C:232:TYR:O	2.07	0.54
1:C:19:PRO:HB3	1:C:79:GLU:HB3	1.88	0.54
1:A:58:THR:OG1	1:A:130:PHE:HB2	2.08	0.54
1:A:26:LEU:HD22	1:A:132:ILE:HD11	1.88	0.54
1:C:114:ALA:HB1	1:C:160:PHE:CE1	2.44	0.53
2:B:231:GLY:O	2:B:233:GLU:N	2.41	0.53
1:A:486:LEU:HD23	1:A:495:ILE:HD11	1.89	0.53
1:A:101:LYS:N	1:A:101:LYS:HD3	2.23	0.53
2:B:312:GLU:HB3	2:B:313:PRO:HD2	1.89	0.53
3:T:716:DA:H2'	3:T:717:DC:C6	2.43	0.53
1:A:253:THR:HG23	1:A:256:ASP:H	1.74	0.53
2:B:35:VAL:O	2:B:39:THR:HG23	2.07	0.53
2:D:191:SER:OG	2:D:198:HIS:ND1	2.23	0.53
2:D:195:ILE:O	2:D:199:ARG:HG3	2.08	0.53
2:D:423:VAL:CG1	2:D:426:TRP:CD1	2.90	0.53
1:A:242:GLN:HB3	1:A:243:PRO:HD2	1.90	0.53
2:B:118:VAL:HG12	2:B:119:PRO:O	2.08	0.53
1:A:433:PRO:HD3	1:A:532:TYR:CZ	2.44	0.53
1:A:326:ILE:O	1:A:341:ILE:HA	2.08	0.53
1:A:5:ILE:HG12	1:A:119:PRO:CD	2.38	0.53
1:C:34:LEU:HD13	1:C:73:LYS:NZ	2.24	0.53
1:A:438:GLU:HB3	1:A:440:PHE:HE2	1.73	0.53
4:P:803:DC:H3'	4:P:804:DA:O4'	2.09	0.53
3:T:706:DA:H2	4:P:822:ATM:HN3	1.48	0.53
2:B:175:ASN:HD21	2:B:201:LYS:NZ	2.07	0.53
2:B:191:SER:OG	2:B:198:HIS:ND1	2.37	0.53
2:D:167:ILE:HD11	2:D:209:LEU:HA	1.91	0.53
1:A:18:GLY:HA3	1:A:56:TYR:CE1	2.43	0.53
1:C:94:ILE:CG2	1:C:95:PRO:HD2	2.39	0.53
2:D:332:GLN:NE2	2:D:427:TYR:HB3	2.24	0.53
1:C:24:TRP:CD2	1:C:25:PRO:HD2	2.43	0.53
2:B:13:LYS:HG3	2:B:16:MET:CE	2.38	0.52
1:C:253:THR:HG22	1:C:256:ASP:CG	2.30	0.52
2:D:46:LYS:HE2	2:D:116:PHE:HB3	1.91	0.52
1:A:229:TRP:CE3	5:A:901:NVP:HCD2	2.44	0.52
2:B:88:TRP:CE3	2:B:89:GLU:HG3	2.44	0.52
2:B:12:LEU:CD1	2:B:124:PHE:HE1	2.23	0.52
1:C:454:LYS:HE2	1:C:554:ALA:O	2.10	0.52
1:C:56:TYR:HB2	1:C:129:ALA:HB3	1.92	0.52
1:A:476:LYS:O	1:A:480:GLN:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:358:ARG:HG3	2:B:366:LYS:HD3	1.92	0.52
1:C:442:VAL:HB	1:C:481:ALA:HB1	1.91	0.52
1:A:85:GLN:C	1:A:154:LYS:HZ1	2.13	0.52
1:A:391:LEU:HD12	1:A:414:TRP:CE3	2.45	0.52
1:C:58:THR:O	1:C:130:PHE:HB2	2.09	0.52
1:C:181:TYR:CE2	1:C:183:TYR:HB2	2.45	0.52
2:B:160:PHE:CD2	2:B:160:PHE:O	2.63	0.52
2:B:231:GLY:C	2:B:233:GLU:N	2.63	0.52
2:B:281:LYS:O	2:B:284:ARG:HG2	2.10	0.52
1:A:195:ILE:O	1:A:199:ARG:NH1	2.43	0.52
1:C:195:ILE:HD11	1:C:199:ARG:HH21	1.75	0.52
2:D:244:ILE:HB	2:D:310:LEU:HD22	1.91	0.52
1:C:65:LYS:HG2	1:C:66:LYS:N	2.25	0.52
1:A:5:ILE:HD11	1:A:119:PRO:HD2	1.92	0.51
1:A:148:VAL:O	1:A:150:PRO:HD3	2.11	0.51
1:A:26:LEU:HD13	1:A:60:VAL:HG12	1.92	0.51
1:C:91:GLN:OE1	1:C:161:GLN:NE2	2.43	0.51
1:C:328:GLU:HG3	1:C:390:LYS:CB	2.39	0.51
3:T:718:DA:H2"	3:T:719:DG:OP2	2.09	0.51
2:D:168:LEU:HD13	2:D:172:LYS:HE3	1.91	0.51
1:A:21:VAL:H	1:A:57:ASN:HB3	1.75	0.51
1:A:179:VAL:HG12	5:A:901:NVP:HCB1	1.91	0.51
1:A:424:LYS:HE2	1:A:426:TRP:CZ3	2.45	0.51
2:B:345:PRO:HB2	2:B:346:PHE:CD2	2.45	0.51
1:C:339:TYR:CZ	1:C:352:GLY:HA3	2.46	0.51
2:D:421:PRO:CB	2:D:423:VAL:H	2.18	0.51
1:C:28:GLU:O	1:C:31:ILE:HG12	2.10	0.51
1:A:102:LYS:HG2	1:A:236:PRO:O	2.11	0.51
2:B:252:TRP:CZ3	2:B:260:LEU:HD22	2.44	0.51
2:D:109:LEU:HD22	2:D:216:THR:HG21	1.92	0.51
1:C:406:TRP:CH2	1:C:407:GLN:HG3	2.43	0.51
1:A:257:ILE:O	1:A:260:LEU:HB3	2.10	0.51
1:C:203:GLU:OE1	1:C:206:ARG:HD2	2.11	0.51
1:A:136:ASN:O	1:A:137:ASN:HB2	2.10	0.51
2:D:78:ARG:HD3	2:D:411:ILE:O	2.11	0.51
1:C:144:TYR:N	1:C:144:TYR:CD2	2.78	0.51
1:C:10:VAL:O	1:C:11:LYS:HG3	2.10	0.51
1:C:34:LEU:CD2	1:C:73:LYS:HD2	2.41	0.51
1:C:209:LEU:O	1:C:214:LEU:HB2	2.11	0.51
1:C:169:GLU:N	1:C:170:PRO:HD2	2.26	0.51
2:B:13:LYS:NZ	2:B:85:GLN:HB3	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:PHE:CD2	1:A:160:PHE:C	2.83	0.51
1:C:91:GLN:HA	1:C:91:GLN:NE2	2.25	0.51
1:C:221:HIS:HB3	1:C:224:GLU:O	2.10	0.50
1:A:447:ASN:OD1	1:A:450:THR:HG23	2.11	0.50
2:D:317:VAL:HG12	2:D:347:LYS:HB3	1.92	0.50
2:B:207:GLN:O	2:B:210:LEU:HB3	2.11	0.50
4:P:816:DG:C2'	4:P:817:MRG:H5'	2.38	0.50
2:B:120:LEU:HD12	2:B:149:LEU:HD23	1.93	0.50
3:E:719:DG:C4'	3:E:720:DG:OP1	2.57	0.50
1:A:221:HIS:HD2	1:A:227:PHE:HE1	1.60	0.50
1:A:160:PHE:HE2	1:A:164:MET:HE2	1.76	0.50
2:D:190:GLY:HA2	2:D:198:HIS:HE1	1.75	0.50
1:A:130:PHE:HD2	1:A:144:TYR:CE2	2.30	0.50
2:D:235:HIS:N	2:D:236:PRO:HD3	2.26	0.50
1:A:329:ILE:O	1:A:392:PRO:HD3	2.12	0.50
1:C:121:ASP:O	1:C:122:GLU:C	2.49	0.50
2:D:168:LEU:HD13	2:D:172:LYS:CE	2.42	0.50
2:D:290:THR:CG2	2:D:290:THR:O	2.60	0.50
4:P:807:DC:H2''	4:P:808:DC:H5'	1.92	0.50
1:A:379:SER:CB	1:A:387:PRO:HD3	2.42	0.50
1:A:144:TYR:CD2	1:A:144:TYR:N	2.77	0.50
1:C:115:TYR:CE1	1:C:160:PHE:HD1	2.30	0.50
4:F:815:DG:H2'	4:F:815:DG:OP2	2.12	0.50
4:P:811:DG:H2'	4:P:812:DT:H72	1.94	0.50
1:A:417:VAL:O	1:A:417:VAL:HG13	2.12	0.49
2:B:171:PHE:HE1	2:B:175:ASN:OD1	1.95	0.49
1:A:5:ILE:HD11	1:A:118:VAL:HA	1.95	0.49
1:C:2:ILE:N	1:C:2:ILE:CD1	2.61	0.49
4:P:803:DC:H2'	4:P:804:DA:C8	2.47	0.49
2:B:423:VAL:HG12	2:B:423:VAL:O	2.11	0.49
2:B:252:TRP:HB3	2:B:257:ILE:HD11	1.94	0.49
1:A:220:LYS:HG3	1:A:220:LYS:O	2.11	0.49
1:A:88:TRP:CD1	2:B:143:ARG:CZ	2.95	0.49
1:C:473:THR:HG23	1:C:476:LYS:HE3	1.94	0.49
1:C:420:PRO:HA	1:C:421:PRO:C	2.33	0.49
1:C:305:GLU:O	1:C:309:ILE:HG23	2.11	0.49
2:B:235:HIS:N	2:B:236:PRO:HD3	2.28	0.49
3:T:715:DA:OP2	3:T:715:DA:H2'	2.13	0.49
1:A:144:TYR:HD2	1:A:144:TYR:N	2.10	0.49
2:D:106:VAL:O	2:D:233:GLU:HA	2.13	0.49
2:B:115:TYR:O	2:B:149:LEU:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:261:VAL:HG22	1:C:276:VAL:CG2	2.43	0.49
1:A:108:VAL:HG12	1:A:227:PHE:CE1	2.47	0.49
2:D:353:LYS:HZ3	2:D:427:TYR:HE1	1.60	0.49
1:A:503:LEU:O	1:A:507:GLN:HB2	2.13	0.49
1:A:279:LEU:HD23	1:A:299:ALA:HB1	1.95	0.49
1:A:317:VAL:HG23	1:A:318:TYR:N	2.28	0.48
2:B:88:TRP:CZ2	2:B:154:LYS:CD	2.96	0.48
2:B:164:MET:CE	2:B:168:LEU:HD11	2.38	0.48
1:C:178:ILE:N	1:C:178:ILE:HD12	2.27	0.48
1:C:164:MET:CE	1:C:168:LEU:HD11	2.42	0.48
1:A:410:TRP:CH2	1:A:412:PRO:HA	2.48	0.48
1:C:279:LEU:HG	1:C:302:GLU:OE1	2.14	0.48
1:C:252:TRP:O	1:C:292:VAL:HG23	2.13	0.48
1:A:328:GLU:HG3	1:A:390:LYS:CB	2.42	0.48
1:C:398:TRP:CZ3	1:C:411:ILE:HD11	2.48	0.48
1:C:184:MET:HE3	4:F:822:ATM:H2'	1.95	0.48
1:C:10:VAL:HG12	1:C:11:LYS:H	1.78	0.48
2:D:279:LEU:HD12	2:D:302:GLU:OE1	2.14	0.48
2:B:286:THR:HG22	2:B:286:THR:O	2.11	0.48
1:C:160:PHE:O	1:C:160:PHE:HD2	1.96	0.48
1:C:253:THR:HG23	1:C:256:ASP:H	1.77	0.48
1:A:424:LYS:HE2	1:A:426:TRP:CE3	2.49	0.48
2:B:425:LEU:CD2	2:B:425:LEU:N	2.76	0.48
1:C:24:TRP:CD2	1:C:25:PRO:CD	2.96	0.48
1:C:202:ILE:O	1:C:206:ARG:HB2	2.14	0.48
1:C:229:TRP:CE3	5:C:901:NVP:HCD2	2.48	0.48
1:C:5:ILE:HD11	1:C:118:VAL:HG22	1.95	0.48
1:C:60:VAL:CG2	1:C:132:ILE:HG21	2.43	0.48
1:C:135:ILE:HG23	1:C:136:ASN:N	2.29	0.48
1:C:331:LYS:HE3	1:C:364:ASP:OD2	2.13	0.48
1:A:540:LYS:HA	1:A:540:LYS:HD3	1.67	0.48
2:B:85:GLN:HA	2:B:88:TRP:NE1	2.29	0.48
1:A:406:TRP:CE3	1:A:407:GLN:HG3	2.46	0.48
2:D:85:GLN:HA	2:D:88:TRP:NE1	2.29	0.48
1:C:130:PHE:CD2	1:C:130:PHE:C	2.87	0.48
1:C:195:ILE:HD11	1:C:199:ARG:NH2	2.28	0.48
1:C:301:LEU:O	1:C:305:GLU:HG2	2.13	0.48
2:B:58:THR:HG21	2:B:75:VAL:HG12	1.95	0.48
1:C:114:ALA:HB1	1:C:160:PHE:HE1	1.79	0.48
1:A:86:ASP:HA	1:A:154:LYS:NZ	2.29	0.48
1:C:10:VAL:HG12	1:C:11:LYS:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:18:GLY:HA3	2:D:127:TYR:CD1	2.48	0.48
2:D:38:CYS:SG	2:D:132:ILE:HD11	2.54	0.48
1:A:115:TYR:CE1	1:A:160:PHE:HD1	2.31	0.47
2:D:167:ILE:HD12	2:D:212:TRP:HB2	1.96	0.47
2:D:395:LYS:HG3	2:D:416:PHE:CE2	2.49	0.47
2:D:105:SER:OG	2:D:235:HIS:CE1	2.67	0.47
2:D:275:LYS:HD2	2:D:277:ARG:NE	2.30	0.47
2:D:363:ASN:O	2:D:367:GLN:HG3	2.14	0.47
1:C:129:ALA:HB1	1:C:143:ARG:HD2	1.96	0.47
1:C:60:VAL:CG2	1:C:130:PHE:HE1	2.26	0.47
2:B:423:VAL:C	2:B:425:LEU:H	2.17	0.47
1:A:34:LEU:HD21	1:A:132:ILE:HD13	1.97	0.47
1:C:195:ILE:O	1:C:199:ARG:HG3	2.14	0.47
2:D:46:LYS:HG2	2:D:148:VAL:HG21	1.96	0.47
1:C:160:PHE:CE2	1:C:164:MET:HB2	2.50	0.47
1:C:37:ILE:HA	1:C:40:GLU:HG3	1.95	0.47
1:A:199:ARG:NE	1:A:222:GLN:OE1	2.38	0.47
2:D:168:LEU:CB	2:D:172:LYS:HE3	2.36	0.47
2:D:423:VAL:HG13	2:D:426:TRP:HE1	1.76	0.47
2:B:421:PRO:HB3	2:B:424:LYS:HB2	1.97	0.47
2:B:207:GLN:HA	2:B:210:LEU:HB3	1.96	0.47
1:A:296:THR:HG23	1:A:299:ALA:H	1.79	0.47
2:B:116:PHE:O	2:B:148:VAL:HG11	2.14	0.47
1:A:230:MET:CE	4:P:821:DC:H1'	2.45	0.47
1:C:497:THR:O	1:C:535:TRP:HA	2.14	0.47
1:C:144:TYR:N	1:C:144:TYR:HD2	2.13	0.47
2:B:142:ILE:CG2	2:B:144:TYR:CE2	2.94	0.47
1:A:197:GLN:O	1:A:200:THR:HB	2.15	0.47
1:C:238:LYS:HD2	1:C:315:HIS:ND1	2.30	0.47
1:C:93:GLY:C	2:D:137:ASN:HD22	2.18	0.47
1:C:296:THR:HG22	1:C:298:GLU:H	1.80	0.47
1:A:314:VAL:HG22	1:A:315:HIS:N	2.30	0.47
1:A:88:TRP:CZ2	2:B:57:ASN:CB	2.98	0.47
2:B:58:THR:CG2	2:B:75:VAL:HG12	2.45	0.47
1:A:230:MET:HE1	4:P:821:DC:H1'	1.96	0.47
1:A:277:ARG:NH1	1:A:334:GLN:O	2.48	0.47
2:D:207:GLN:OE1	2:D:210:LEU:HD23	2.14	0.47
1:A:130:PHE:O	1:A:143:ARG:HG2	2.14	0.47
1:C:144:TYR:HD2	1:C:144:TYR:H	1.63	0.47
3:E:717:DC:H2'	3:E:718:DA:H8	1.75	0.47
1:A:365:VAL:HG11	1:A:401:TRP:CG	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:101:LYS:HD3	1:C:101:LYS:N	2.30	0.47
2:D:242:GLN:O	2:D:242:GLN:HG3	2.14	0.47
2:B:425:LEU:HD13	2:B:425:LEU:HA	1.46	0.47
2:B:58:THR:HG23	2:B:76:ASP:O	2.15	0.47
1:C:317:VAL:HG23	1:C:349:LEU:HD23	1.97	0.47
1:C:65:LYS:HD3	1:C:72:ARG:HH22	1.80	0.46
2:B:131:THR:HG23	2:B:141:GLY:HA3	1.97	0.46
1:A:184:MET:HG3	4:P:822:ATM:N4'	2.30	0.46
1:C:171:PHE:CD1	1:C:205:LEU:HD12	2.50	0.46
3:E:716:DA:H61	4:F:812:DT:H3	1.64	0.46
2:B:214:LEU:HD12	2:B:215:THR:N	2.29	0.46
1:C:120:LEU:CG	1:C:121:ASP:N	2.76	0.46
1:A:149:LEU:HA	1:A:150:PRO:HD3	1.72	0.46
2:B:28:GLU:CG	2:B:135:ILE:HD11	2.46	0.46
4:F:807:DC:H2''	4:F:808:DC:C5'	2.45	0.46
1:C:406:TRP:CZ3	1:C:407:GLN:HG3	2.51	0.46
1:A:135:ILE:HG23	1:A:136:ASN:N	2.18	0.46
2:B:115:TYR:HE1	2:B:157:PRO:N	2.13	0.46
2:B:32:LYS:HD3	2:B:32:LYS:HA	1.55	0.46
2:D:350:LYS:HE3	2:D:378:GLU:OE2	2.15	0.46
1:C:171:PHE:CD2	1:C:171:PHE:C	2.89	0.46
1:A:382:ILE:O	2:B:136:ASN:HB2	2.16	0.46
4:F:816:DG:C2'	4:F:817:MRG:H5'	2.36	0.46
2:D:209:LEU:HD22	2:D:214:LEU:HD23	1.98	0.46
1:C:75:VAL:HG12	1:C:77:PHE:CD1	2.44	0.46
2:B:209:LEU:H	2:B:209:LEU:HD12	1.79	0.46
2:B:427:TYR:O	2:B:428:GLN:O	2.34	0.46
1:C:47:ILE:HD13	1:C:144:TYR:CE1	2.50	0.46
1:A:115:TYR:HD1	1:A:160:PHE:HD1	1.61	0.46
1:C:274:ILE:CD1	1:C:309:ILE:HD11	2.44	0.46
2:D:319:TYR:CD2	2:D:383:TRP:HD1	2.34	0.46
1:A:455:ALA:CB	1:A:469:LEU:HD11	2.45	0.46
1:A:442:VAL:HB	1:A:481:ALA:HB1	1.98	0.46
2:D:381:VAL:O	2:D:381:VAL:HG12	2.15	0.46
1:C:34:LEU:HD22	1:C:73:LYS:HD2	1.97	0.46
2:B:34:LEU:HA	2:B:34:LEU:HD23	1.83	0.46
1:C:307:ARG:O	1:C:311:LYS:HG3	2.16	0.45
2:D:57:ASN:HA	2:D:129:ALA:O	2.16	0.45
2:D:298:GLU:CD	2:D:298:GLU:H	2.20	0.45
2:D:84:THR:O	2:D:84:THR:HG22	2.16	0.45
1:A:134:SER:OG	1:A:135:ILE:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:195:ILE:HG12	2:B:199:ARG:HE	1.81	0.45
3:T:716:DA:C2'	3:T:717:DC:H6	2.28	0.45
2:B:103:LYS:HE2	2:B:179:VAL:HG23	1.98	0.45
1:A:8:VAL:O	1:A:10:VAL:HG23	2.16	0.45
1:C:296:THR:HG22	1:C:297:GLU:N	2.31	0.45
1:C:47:ILE:HD13	1:C:144:TYR:CD1	2.52	0.45
1:C:24:TRP:HA	1:C:25:PRO:HD3	1.74	0.45
1:A:254:VAL:HG23	1:A:289:LEU:HD12	1.98	0.45
2:D:423:VAL:O	2:D:426:TRP:HD1	1.99	0.45
1:A:497:THR:O	1:A:535:TRP:HA	2.16	0.45
1:A:171:PHE:CD2	1:A:171:PHE:C	2.89	0.45
2:B:328:GLU:O	2:B:339:TYR:HA	2.16	0.45
2:D:85:GLN:HA	2:D:88:TRP:CE2	2.51	0.45
1:C:159:ILE:O	1:C:163:SER:OG	2.34	0.45
2:B:87:PHE:CD2	2:B:87:PHE:C	2.89	0.45
2:B:424:LYS:C	2:B:425:LEU:HD22	2.36	0.45
2:B:382:ILE:HG22	2:B:383:TRP:CE2	2.52	0.45
1:A:109:LEU:HD21	1:A:202:ILE:HG23	1.97	0.45
2:D:115:TYR:CE1	2:D:156:SER:C	2.90	0.45
2:D:115:TYR:O	2:D:149:LEU:HB2	2.17	0.45
1:C:254:VAL:HG21	1:C:286:THR:HG21	1.97	0.45
1:C:5:ILE:HD11	1:C:119:PRO:HD2	1.98	0.45
1:C:296:THR:HG22	1:C:298:GLU:HG2	1.99	0.45
4:P:812:DT:H2'	4:P:813:DT:H71	1.98	0.45
1:A:324:ASP:O	1:A:343:GLN:HG2	2.16	0.45
2:D:356:ARG:HG3	2:D:356:ARG:O	2.17	0.44
2:B:13:LYS:HD3	2:B:86:ASP:H	1.82	0.44
2:B:317:VAL:O	2:B:317:VAL:HG23	2.16	0.44
2:D:252:TRP:CD1	2:D:295:LEU:CD1	3.01	0.44
2:D:421:PRO:HB2	2:D:423:VAL:HB	1.99	0.44
1:A:160:PHE:CE2	1:A:164:MET:CE	3.00	0.44
1:A:47:ILE:CG2	1:A:48:SER:N	2.80	0.44
1:A:108:VAL:HG13	1:A:108:VAL:O	2.18	0.44
1:C:55:PRO:O	1:C:56:TYR:CG	2.70	0.44
1:A:204:GLU:O	1:A:207:GLN:HB2	2.18	0.44
1:A:104:LYS:HB2	1:A:104:LYS:HE3	1.62	0.44
2:D:266:TRP:CD1	2:D:422:LEU:HD21	2.53	0.44
2:D:114:ALA:H	2:D:214:LEU:HD13	1.81	0.44
1:C:24:TRP:CG	1:C:25:PRO:HD2	2.52	0.44
2:D:78:ARG:O	2:D:82:LYS:HG3	2.17	0.44
1:C:130:PHE:CD2	1:C:144:TYR:HE2	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLU:N	1:A:170:PRO:HD2	2.32	0.44
2:B:252:TRP:CB	2:B:257:ILE:HD11	2.48	0.44
1:C:120:LEU:CD1	1:C:121:ASP:H	2.30	0.44
1:C:440:PHE:CZ	1:C:489:SER:HB3	2.53	0.44
1:C:49:LYS:C	1:C:50:ILE:HD12	2.38	0.44
2:D:357:MET:O	2:D:360:ALA:HB2	2.18	0.44
1:A:253:THR:HG23	1:A:256:ASP:N	2.32	0.44
1:C:21:VAL:O	1:C:22:LYS:HD3	2.17	0.44
1:A:26:LEU:HD23	1:A:31:ILE:HG22	1.99	0.44
1:C:12:LEU:HD11	1:C:127:TYR:CE1	2.53	0.44
2:D:101:LYS:O	2:D:236:PRO:CB	2.66	0.44
1:A:253:THR:HG22	1:A:256:ASP:CG	2.38	0.44
2:B:210:LEU:HD23	2:B:211:ARG:HG3	1.99	0.44
1:C:21:VAL:HG12	1:C:22:LYS:N	2.23	0.44
2:B:209:LEU:CD2	2:B:214:LEU:HD23	2.45	0.44
2:B:423:VAL:HG13	2:B:426:TRP:HE1	1.81	0.44
1:C:78:ARG:HB3	1:C:82:LYS:NZ	2.33	0.44
1:C:320:ASP:HB3	1:C:323:LYS:HE3	1.99	0.44
1:A:32:LYS:CD	1:A:135:ILE:HB	2.39	0.44
1:A:252:TRP:CE2	1:A:295:LEU:HD11	2.53	0.44
2:D:11:LYS:HG2	2:D:12:LEU:O	2.18	0.44
1:A:234:LEU:HD23	1:A:318:TYR:OH	2.18	0.44
2:D:422:LEU:O	2:D:422:LEU:HG	2.18	0.43
1:A:160:PHE:CE2	1:A:164:MET:HE2	2.53	0.43
2:D:323:LYS:HD2	2:D:344:GLU:OE2	2.18	0.43
2:D:353:LYS:NZ	2:D:427:TYR:HE1	2.15	0.43
2:D:167:ILE:HD13	2:D:209:LEU:HD23	2.00	0.43
4:F:807:DC:H2'	4:F:808:DC:C6	2.53	0.43
1:C:434:ILE:CD1	1:C:530:LYS:HB3	2.47	0.43
1:A:199:ARG:HE	1:A:222:GLN:CD	2.17	0.43
2:B:88:TRP:CH2	2:B:154:LYS:HD3	2.53	0.43
2:D:423:VAL:CG1	2:D:426:TRP:HD1	2.31	0.43
1:A:130:PHE:CZ	1:A:132:ILE:HB	2.53	0.43
2:B:421:PRO:HB3	2:B:424:LYS:HD2	2.00	0.43
2:B:234:LEU:C	2:B:236:PRO:HD3	2.39	0.43
2:B:345:PRO:HB2	2:B:346:PHE:HD2	1.83	0.43
1:C:334:GLN:O	1:C:336:GLN:HG3	2.18	0.43
1:A:108:VAL:CG1	1:A:227:PHE:CE1	3.02	0.43
2:D:21:VAL:HG12	2:D:59:PRO:HG3	1.99	0.43
1:A:451:LYS:O	1:A:471:ASN:N	2.51	0.43
1:C:100:LEU:O	1:C:318:TYR:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:17:ASP:OD1	1:C:127:TYR:HE1	2.02	0.43
2:D:320:ASP:OD1	2:D:322:SER:CB	2.66	0.43
1:C:249:LYS:O	1:C:252:TRP:CE2	2.72	0.43
1:C:37:ILE:HG23	1:C:40:GLU:OE1	2.18	0.43
2:B:103:LYS:HE2	2:B:179:VAL:CG2	2.48	0.43
3:T:706:DA:C6	4:P:822:ATM:O4	2.71	0.43
1:C:225:PRO:CB	1:C:226:PRO:HA	2.48	0.43
2:D:81:ASN:OD1	2:D:153:TRP:HD1	2.01	0.43
1:C:24:TRP:CG	1:C:25:PRO:CD	3.02	0.43
1:A:328:GLU:HG3	1:A:390:LYS:HB3	2.01	0.43
1:C:246:LEU:HD22	1:C:260:LEU:CD1	2.49	0.43
1:C:194:GLU:HB2	1:C:197:GLN:HG3	2.00	0.43
2:D:34:LEU:HD23	2:D:34:LEU:HA	1.83	0.43
2:D:169:GLU:HB3	2:D:170:PRO:HD3	2.01	0.43
2:D:358:ARG:CG	2:D:366:LYS:HD3	2.47	0.43
1:C:164:MET:CE	1:C:187:LEU:HD11	2.48	0.43
2:B:336:GLN:NE2	2:B:427:TYR:CE1	2.87	0.43
2:D:328:GLU:HG3	2:D:342:TYR:HE2	1.83	0.43
2:D:422:LEU:HD23	2:D:423:VAL:N	2.33	0.43
1:C:60:VAL:HG23	1:C:130:PHE:CE1	2.47	0.43
1:C:271:TYR:CD1	1:C:314:VAL:CG1	3.02	0.43
2:D:362:THR:HG23	2:D:363:ASN:H	1.84	0.43
1:C:88:TRP:CE3	1:C:88:TRP:HA	2.53	0.43
1:A:178:ILE:HG22	1:A:179:VAL:N	2.34	0.43
2:B:203:GLU:O	2:B:207:GLN:HG2	2.18	0.43
2:B:167:ILE:HD13	2:B:212:TRP:HB3	2.01	0.43
1:A:253:THR:O	1:A:256:ASP:N	2.51	0.43
1:C:339:TYR:CE2	1:C:352:GLY:HA3	2.54	0.43
4:P:816:DG:H2'	4:P:817:MRG:H8	2.01	0.42
2:B:114:ALA:CA	2:B:214:LEU:HD13	2.48	0.42
2:D:382:ILE:HG22	2:D:383:TRP:CE2	2.54	0.42
1:A:132:ILE:CG1	1:A:133:PRO:HD2	2.48	0.42
3:T:714:DG:H2''	3:T:715:DA:C8	2.54	0.42
2:B:78:ARG:HD3	2:B:411:ILE:O	2.18	0.42
3:T:706:DA:H2'	3:T:707:DG:C8	2.54	0.42
4:F:807:DC:H4'	4:F:808:DC:OP1	2.19	0.42
2:D:355:ALA:O	2:D:357:MET:N	2.52	0.42
1:C:76:ASP:O	1:C:76:ASP:OD1	2.37	0.42
1:C:246:LEU:HA	1:C:247:PRO:HD3	1.81	0.42
1:C:102:LYS:HG3	1:C:237:ASP:HA	2.00	0.42
1:C:447:ASN:HD21	1:C:450:THR:HG23	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:SER:OG	1:A:137:ASN:HA	2.18	0.42
1:C:184:MET:HG3	4:F:822:ATM:N4'	2.34	0.42
1:C:84:THR:HG22	1:C:85:GLN:N	2.34	0.42
1:C:35:VAL:HG23	1:C:134:SER:OG	2.19	0.42
2:D:387:PRO:HG2	2:D:389:PHE:CE2	2.54	0.42
1:A:118:VAL:HB	1:A:149:LEU:CD1	2.48	0.42
3:E:721:DG:H2''	3:E:722:DA:C8	2.54	0.42
1:C:184:MET:HE1	4:F:822:ATM:H2'	2.01	0.42
2:B:297:GLU:O	2:B:301:LEU:N	2.48	0.42
2:B:279:LEU:HD11	2:B:303:LEU:HB2	2.02	0.42
2:D:292:VAL:C	2:D:293:ILE:HD13	2.39	0.42
1:A:473:THR:HG23	1:A:476:LYS:HG3	2.01	0.42
1:C:448:ARG:O	1:C:451:LYS:HD3	2.19	0.42
2:B:61:PHE:CD1	2:B:403:THR:HG22	2.55	0.42
1:A:418:ASN:OD1	1:A:420:PRO:HD3	2.19	0.42
2:D:345:PRO:O	2:D:346:PHE:HB2	2.18	0.42
2:B:65:LYS:O	2:B:68:SER:HB3	2.20	0.42
1:C:540:LYS:HA	1:C:540:LYS:HD2	1.87	0.42
2:B:115:TYR:N	2:B:115:TYR:CD2	2.88	0.42
2:D:423:VAL:HG12	2:D:423:VAL:O	2.18	0.42
1:C:160:PHE:HE2	1:C:164:MET:HB2	1.84	0.42
1:C:104:LYS:HB2	1:C:104:LYS:HE3	1.81	0.42
1:C:21:VAL:HG23	1:C:57:ASN:O	2.19	0.42
2:B:28:GLU:HG3	2:B:32:LYS:HE2	2.02	0.42
1:A:512:LYS:HG2	1:A:513:SER:N	2.34	0.42
2:D:12:LEU:HD12	2:D:124:PHE:HE1	1.84	0.42
2:B:336:GLN:NE2	2:B:427:TYR:CZ	2.88	0.42
1:C:94:ILE:HG22	1:C:95:PRO:HD2	2.01	0.42
1:A:328:GLU:HG3	1:A:390:LYS:HB2	2.01	0.42
1:C:354:TYR:HD1	1:C:374:LYS:HD2	1.84	0.42
1:C:172:LYS:HE2	1:C:180:ILE:HD12	2.02	0.42
2:D:209:LEU:HD22	2:D:214:LEU:CD2	2.50	0.42
2:D:320:ASP:OD2	2:D:323:LYS:CE	2.68	0.42
1:C:164:MET:HE2	1:C:168:LEU:CD1	2.50	0.42
1:A:394:GLN:NE2	1:A:416:PHE:CE2	2.88	0.42
1:C:37:ILE:HG12	1:C:40:GLU:OE2	2.20	0.42
1:C:37:ILE:HG22	1:C:37:ILE:O	2.19	0.42
2:B:154:LYS:O	2:B:157:PRO:HD2	2.20	0.41
1:A:48:SER:O	1:A:144:TYR:HB2	2.20	0.41
1:C:424:LYS:HE2	1:C:426:TRP:CE3	2.55	0.41
2:D:160:PHE:O	2:D:160:PHE:CG	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:MET:HE2	1:A:512:LYS:HD3	2.02	0.41
1:A:105:SER:O	1:A:190:GLY:HA2	2.20	0.41
2:D:56:TYR:O	2:D:143:ARG:NH2	2.53	0.41
2:B:380:ILE:O	2:B:384:GLY:HA2	2.19	0.41
2:D:287:LYS:HE3	2:D:291:GLU:OE2	2.20	0.41
1:C:23:GLN:NE2	1:C:133:PRO:HD2	2.34	0.41
3:E:716:DA:H2''	3:E:717:DC:OP2	2.20	0.41
2:B:331:LYS:O	2:B:424:LYS:HE3	2.21	0.41
1:C:164:MET:CE	1:C:168:LEU:CD1	2.98	0.41
1:C:34:LEU:HD13	1:C:73:LYS:HZ1	1.84	0.41
2:D:5:ILE:H	2:D:5:ILE:HG13	1.66	0.41
1:A:130:PHE:CG	1:A:131:THR:N	2.88	0.41
1:C:39:THR:HA	1:C:42:GLU:OE1	2.20	0.41
1:A:257:ILE:CG2	1:A:283:LEU:HD21	2.51	0.41
1:A:77:PHE:O	1:A:78:ARG:C	2.58	0.41
1:C:270:ILE:HG23	1:C:271:TYR:N	2.36	0.41
1:A:450:THR:OG1	1:A:452:LEU:HB2	2.20	0.41
4:F:817:MRG:H2''	4:F:818:DC:O5'	2.20	0.41
2:D:115:TYR:OH	2:D:157:PRO:HB3	2.21	0.41
2:B:301:LEU:O	2:B:305:GLU:HB2	2.21	0.41
1:C:65:LYS:HB3	1:C:72:ARG:NH1	2.36	0.41
2:B:344:GLU:O	2:B:345:PRO:C	2.59	0.41
2:B:360:ALA:O	2:B:361:HIS:HB2	2.20	0.41
2:B:370:GLU:HA	2:B:370:GLU:OE1	2.20	0.41
2:D:355:ALA:O	2:D:356:ARG:C	2.59	0.41
2:D:201:LYS:HB2	2:D:201:LYS:HE3	1.90	0.41
2:D:423:VAL:O	2:D:426:TRP:CD1	2.74	0.41
2:D:167:ILE:HG12	2:D:167:ILE:O	2.19	0.41
1:A:34:LEU:HD22	1:A:73:LYS:CE	2.41	0.41
1:C:26:LEU:HD22	1:C:60:VAL:HG12	2.03	0.41
1:C:457:TYR:C	1:C:457:TYR:HD1	2.24	0.41
1:A:51:GLY:N	1:A:52:PRO:HD2	2.36	0.41
2:B:353:LYS:HZ2	2:B:427:TYR:HE1	1.69	0.41
2:B:345:PRO:O	2:B:346:PHE:HB2	2.21	0.41
1:A:416:PHE:CD1	1:A:417:VAL:N	2.89	0.41
1:A:339:TYR:CZ	1:A:352:GLY:HA3	2.56	0.41
1:C:516:GLU:O	1:C:520:GLN:HG3	2.21	0.41
1:C:416:PHE:CZ	1:C:418:ASN:HB2	2.56	0.41
4:F:803:DC:H2'	4:F:803:DC:O2	2.20	0.41
1:A:144:TYR:HD2	1:A:144:TYR:H	1.67	0.41
1:A:28:GLU:HA	1:A:31:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:60:VAL:HG21	1:C:132:ILE:HG21	2.01	0.41
1:C:107:THR:HG23	1:C:222:GLN:O	2.20	0.41
3:E:713:DC:H2''	3:E:714:DG:H5'	2.03	0.41
2:D:207:GLN:O	2:D:210:LEU:N	2.53	0.41
4:F:819:DG:H2'	4:F:820:DC:C6	2.55	0.41
1:C:103:LYS:O	1:C:236:PRO:HB3	2.21	0.41
2:D:50:ILE:HD12	2:D:54:ASN:CB	2.51	0.41
1:C:271:TYR:CE1	1:C:314:VAL:CG1	3.03	0.40
1:C:476:LYS:O	1:C:480:GLN:HB2	2.21	0.40
2:D:50:ILE:HD12	2:D:54:ASN:HB3	2.04	0.40
1:C:22:LYS:O	1:C:23:GLN:HB2	2.20	0.40
1:A:28:GLU:HA	1:A:31:ILE:CD1	2.51	0.40
1:A:398:TRP:CE2	1:A:411:ILE:HD12	2.56	0.40
1:C:424:LYS:HE2	1:C:426:TRP:CZ3	2.57	0.40
1:A:308:GLU:HA	1:A:311:LYS:HG3	2.02	0.40
1:C:276:VAL:HG22	1:C:276:VAL:O	2.21	0.40
1:C:58:THR:CG2	1:C:77:PHE:HB3	2.51	0.40
2:B:423:VAL:O	2:B:423:VAL:CG1	2.68	0.40
1:A:460:ASN:HA	2:B:286:THR:HG22	2.04	0.40
1:C:8:VAL:HA	1:C:9:PRO:HD3	1.97	0.40
1:A:323:LYS:HE3	1:A:344:GLU:HG2	2.03	0.40
1:C:46:LYS:HD3	1:C:116:PHE:HB3	2.03	0.40
2:B:120:LEU:HD23	2:B:125:ARG:CG	2.52	0.40
1:C:226:PRO:HA	1:C:234:LEU:O	2.21	0.40
1:A:221:HIS:CG	1:A:224:GLU:CB	3.02	0.40
1:C:220:LYS:HE2	1:C:222:GLN:CG	2.51	0.40
2:D:275:LYS:HD3	2:D:275:LYS:HA	1.81	0.40
2:B:331:LYS:HB2	2:B:337:TRP:CZ3	2.57	0.40
1:C:74:LEU:HD23	1:C:151:GLN:HB3	2.03	0.40
1:A:15:GLY:O	1:A:16:MET:HG2	2.22	0.40
1:A:33:ALA:O	1:A:37:ILE:HG13	2.22	0.40
1:C:255:ASN:HB2	1:C:289:LEU:HG	2.02	0.40
2:B:115:TYR:HD2	2:B:115:TYR:H	1.69	0.40
1:A:406:TRP:CH2	2:B:418:ASN:CA	3.01	0.40
1:A:54:ASN:HD22	1:A:129:ALA:CB	2.35	0.40
1:A:221:HIS:HB2	1:A:224:GLU:H	1.86	0.40
2:B:57:ASN:HA	2:B:129:ALA:O	2.22	0.40
1:A:21:VAL:HG12	1:A:22:LYS:N	2.36	0.40
1:A:491:LEU:HA	1:A:491:LEU:HD23	1.89	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	553/556 (100%)	524 (95%)	29 (5%)	0	100	100
1	C	553/556 (100%)	522 (94%)	31 (6%)	0	100	100
2	B	408/428 (95%)	395 (97%)	13 (3%)	0	100	100
2	D	408/428 (95%)	396 (97%)	12 (3%)	0	100	100
All	All	1922/1968 (98%)	1837 (96%)	85 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	495/497 (100%)	466 (94%)	29 (6%)	24	54
1	C	495/497 (100%)	471 (95%)	24 (5%)	31	65
2	B	374/390 (96%)	344 (92%)	30 (8%)	15	38
2	D	374/390 (96%)	357 (96%)	17 (4%)	34	67
All	All	1738/1774 (98%)	1638 (94%)	100 (6%)	25	55

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	3	SER

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Mol	Chain	Res	Type
1	A	5	ILE
1	A	58	THR
1	A	107	THR
1	A	110	ASP
1	A	117	SER
1	A	123	ASP
1	A	130	PHE
1	A	137	ASN
1	A	144	TYR
1	A	146	TYR
1	A	159	ILE
1	A	160	PHE
1	A	163	SER
1	A	182	GLN
1	A	205	LEU
1	A	221	HIS
1	A	297	GLU
1	A	329	ILE
1	A	357	MET
1	A	399	GLU
1	A	428	GLN
1	A	459	THR
1	A	463	ARG
1	A	469	LEU
1	A	473	THR
1	A	474	ASN
1	A	539	HIS
2	B	7	THR
2	B	32	LYS
2	B	47	ILE
2	B	50	ILE
2	B	55	PRO
2	B	69	THR
2	B	72	ARG
2	B	87	PHE
2	B	113	ASP
2	B	115	TYR
2	B	123	ASP
2	B	163	SER
2	B	171	PHE
2	B	210	LEU
2	B	232	TYR

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Mol	Chain	Res	Type
2	B	233	GLU
2	B	240	THR
2	B	250	ASP
2	B	274	ILE
2	B	276	VAL
2	B	280	SER
2	B	293	ILE
2	B	305	GLU
2	B	321	PRO
2	B	324	ASP
2	B	336	GLN
2	B	403	THR
2	B	410	TRP
2	B	418	ASN
2	B	427	TYR
1	C	2	ILE
1	C	3	SER
1	C	8	VAL
1	C	101	LYS
1	C	105	SER
1	C	130	PHE
1	C	142	ILE
1	C	144	TYR
1	C	146	TYR
1	C	160	PHE
1	C	163	SER
1	C	182	GLN
1	C	205	LEU
1	C	211	ARG
1	C	215	THR
1	C	270	ILE
1	C	309	ILE
1	C	342	TYR
1	C	357	MET
1	C	428	GLN
1	C	452	LEU
1	C	459	THR
1	C	472	THR
1	C	497	THR
2	D	5	ILE
2	D	163	SER
2	D	194	GLU

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Mol	Chain	Res	Type
2	D	234	LEU
2	D	241	VAL
2	D	248	GLU
2	D	277	ARG
2	D	295	LEU
2	D	296	THR
2	D	317	VAL
2	D	357	MET
2	D	370	GLU
2	D	390	LYS
2	D	417	VAL
2	D	419	THR
2	D	422	LEU
2	D	424	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	221	HIS
1	A	487	GLN
1	A	545	ASN
2	B	175	ASN
2	B	336	GLN
1	C	23	GLN
1	C	175	ASN
1	C	222	GLN
1	C	235	HIS
1	C	447	ASN
2	D	137	ASN
2	D	235	HIS
2	D	269	GLN
2	D	336	GLN
2	D	367	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	MRG	F	817	1,3,4	20,28,29	3.10	9 (45%)	25,39,42	2.03	8 (32%)
4	ATM	F	822	3,4	13,23,24	1.70	2 (15%)	17,32,35	1.85	3 (17%)
4	MRG	P	817	1,3,4	20,28,29	3.17	9 (45%)	25,39,42	2.04	8 (32%)
4	ATM	P	822	3,4	13,23,24	1.56	2 (15%)	17,32,35	1.89	3 (17%)

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
4	MRG	F	817	1,3,4	-	0/8/26/27	0/3/3/3
4	ATM	F	822	3,4	-	0/6/24/25	0/2/2/2
4	MRG	P	817	1,3,4	-	0/8/26/27	0/3/3/3
4	ATM	P	822	3,4	-	0/6/24/25	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	817	MRG	C2'-C3'	-3.09	1.44	1.52
4	F	817	MRG	C2'-C3'	-3.01	1.44	1.52
4	P	817	MRG	O4'-C4'	-2.42	1.39	1.45
4	F	817	MRG	O4'-C4'	-2.39	1.39	1.45
4	F	817	MRG	O3'-C3'	-2.33	1.38	1.43
4	P	817	MRG	O3'-C3'	-2.23	1.38	1.43
4	P	817	MRG	C2-N1	2.37	1.43	1.34
4	F	817	MRG	C2-N1	2.44	1.43	1.34
4	F	817	MRG	C2-N3	2.76	1.44	1.34
4	P	817	MRG	C2-N3	2.94	1.45	1.34
4	P	822	ATM	C4-N3	3.55	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	822	ATM	C6-N1	3.60	1.40	1.35
4	P	817	MRG	C6-C5	3.64	1.48	1.41
4	F	822	ATM	C6-N1	3.77	1.40	1.35
4	F	817	MRG	C6-C5	3.79	1.48	1.41
4	F	822	ATM	C4-N3	3.82	1.40	1.33
4	F	817	MRG	C6-N1	4.65	1.41	1.33
4	P	817	MRG	C6-N1	4.72	1.41	1.33
4	F	817	MRG	C4-N3	6.41	1.45	1.35
4	P	817	MRG	C4-N3	6.84	1.46	1.35
4	F	817	MRG	C2-N2	8.59	1.48	1.34
4	P	817	MRG	C2-N2	8.70	1.48	1.34

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	822	ATM	C5-C4-N3	-4.25	120.40	125.14
4	P	822	ATM	C5-C4-N3	-4.20	120.46	125.14
4	P	817	MRG	N3-C2-N1	-4.15	119.85	126.22
4	F	817	MRG	N3-C2-N1	-4.06	119.98	126.22
4	F	817	MRG	C5-C6-N1	-2.43	120.26	123.59
4	P	817	MRG	C5-C6-N1	-2.35	120.37	123.59
4	F	817	MRG	C4-C5-N7	-2.28	107.38	109.48
4	P	822	ATM	C5'-C4'-C3'	-2.14	106.23	114.46
4	F	822	ATM	C5'-C4'-C3'	-2.07	106.51	114.46
4	P	817	MRG	N2-C2-N3	2.60	122.08	117.22
4	F	817	MRG	C22-C21-N2	2.69	118.74	111.46
4	F	817	MRG	O5'-C5'-C4'	2.89	119.71	109.12
4	P	817	MRG	O5'-C5'-C4'	2.97	120.00	109.12
4	P	817	MRG	C22-C21-N2	3.12	119.92	111.46
4	F	817	MRG	C6-N1-C2	3.19	119.96	115.31
4	P	817	MRG	C6-N1-C2	3.32	120.13	115.31
4	F	817	MRG	O4'-C1'-N9	3.75	114.21	107.72
4	F	817	MRG	C2-N3-C4	4.06	119.98	115.09
4	P	817	MRG	C2-N3-C4	4.10	120.03	115.09
4	P	817	MRG	O4'-C1'-N9	4.16	114.92	107.72
4	F	822	ATM	C4-N3-C2	4.96	119.53	115.25
4	P	822	ATM	C4-N3-C2	5.05	119.61	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	817	MRG	4	0
4	F	822	ATM	5	0
4	P	817	MRG	4	0
4	P	822	ATM	5	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	$\# Z > 2$
5	NVP	A	901	-	18,23,23	1.98	5 (27%)	18,34,34	2.05	7 (38%)
5	NVP	C	901	-	18,23,23	2.00	5 (27%)	18,34,34	2.22	7 (38%)

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
5	NVP	A	901	-	-	0/0/6/6	0/2/4/4
5	NVP	C	901	-	-	0/0/6/6	0/2/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	901	NVP	C10-C15	-2.64	1.38	1.40
5	A	901	NVP	C10-C15	-2.16	1.38	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	901	NVP	C4-N3	3.32	1.39	1.32
5	C	901	NVP	C15-N14	3.52	1.40	1.35
5	A	901	NVP	C4-N3	3.55	1.39	1.32
5	C	901	NVP	C2-N3	3.57	1.40	1.35
5	A	901	NVP	C13-N14	3.72	1.40	1.32
5	A	901	NVP	C15-N14	3.77	1.41	1.35
5	A	901	NVP	C2-N3	3.85	1.41	1.35
5	C	901	NVP	C13-N14	3.94	1.40	1.32

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	901	NVP	CC-CA-N1	-4.02	114.11	118.25
5	A	901	NVP	CC-CA-N1	-3.28	114.87	118.25
5	C	901	NVP	C10-C15-N14	-3.14	119.80	123.53
5	A	901	NVP	CB-CA-N1	-3.12	115.03	118.25
5	A	901	NVP	C7-C2-N3	-2.88	119.75	122.90
5	C	901	NVP	C7-C2-N3	-2.84	119.79	122.90
5	A	901	NVP	C12-C13-N14	-2.51	120.00	123.94
5	A	901	NVP	C10-C15-N14	-2.49	120.57	123.53
5	C	901	NVP	C12-C13-N14	-2.21	120.47	123.94
5	C	901	NVP	CB-CA-N1	-2.19	116.00	118.25
5	C	901	NVP	CD-C6-C7	2.64	122.71	119.84
5	A	901	NVP	CD-C6-C7	3.14	123.25	119.84
5	A	901	NVP	C11-C10-C15	4.07	120.80	117.00
5	C	901	NVP	C11-C10-C15	5.26	121.92	117.00

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
5	A	901	NVP	2	0
5	C	901	NVP	1	0

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	555/556 (99%)	0.77	78 (14%) 4 2	25, 77, 185, 198	0
1	C	555/556 (99%)	0.75	86 (15%) 3 1	24, 79, 184, 196	0
2	B	412/428 (96%)	0.12	18 (4%) 38 31	28, 63, 141, 161	0
2	D	412/428 (96%)	0.17	24 (5%) 26 20	31, 64, 149, 167	0
3	E	20/27 (74%)	0.37	2 (10%) 9 5	78, 107, 166, 175	0
3	T	20/27 (74%)	0.31	3 (15%) 3 2	75, 92, 162, 179	0
4	F	18/21 (85%)	0.48	1 (5%) 28 22	49, 95, 166, 171	0
4	P	18/21 (85%)	0.37	1 (5%) 28 22	49, 88, 164, 175	0
All	All	2010/2064 (97%)	0.49	213 (10%) 8 4	24, 71, 172, 198	0

All (213) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	PRO	20.1
1	A	132	ILE	15.0
1	A	60	VAL	13.7
1	C	62	ALA	13.1
1	C	61	PHE	13.0
1	A	27	THR	12.5
1	C	73	LYS	12.3
1	C	133	PRO	10.4
1	C	136	ASN	10.3
1	A	61	PHE	10.0
1	A	34	LEU	9.6
1	C	132	ILE	9.5
1	A	26	LEU	9.4
1	C	74	LEU	9.0
1	C	31	ILE	8.8
1	A	64	LYS	8.6

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Mol	Chain	Res	Type	RSRZ
1	A	33	ALA	8.5
1	A	67	ASP	8.4
1	A	28	GLU	8.3
1	C	33	ALA	8.3
1	C	64	LYS	8.0
1	A	68	SER	7.9
1	C	34	LEU	7.8
1	A	74	LEU	7.8
1	A	31	ILE	7.7
1	C	72	ARG	7.6
1	A	25	PRO	7.3
1	C	137	ASN	7.1
1	C	59	PRO	6.9
1	C	252	TRP	6.9
1	A	222	GLN	6.9
1	C	27	THR	6.6
1	A	62	ALA	6.6
2	B	214	LEU	6.3
1	C	24	TRP	6.3
1	C	292	VAL	6.2
1	C	65	LYS	6.2
1	C	60	VAL	6.1
1	A	140	PRO	6.0
1	A	139	THR	6.0
1	C	219	LYS	5.9
1	C	223	LYS	5.9
1	A	255	ASN	5.9
1	C	15	GLY	5.8
1	A	32	LYS	5.7
1	C	25	PRO	5.6
1	C	26	LEU	5.5
1	C	71	TRP	5.5
1	A	75	VAL	5.3
2	D	360	ALA	5.3
1	C	32	LYS	5.3
1	A	220	LYS	5.2
1	A	137	ASN	5.2
2	B	93	GLY	5.2
1	A	287	LYS	5.1
1	C	251	SER	5.1
1	A	289	LEU	5.1
1	A	24	TRP	5.0

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Mol	Chain	Res	Type	RSRZ
1	C	69	THR	5.0
2	D	94	ILE	4.9
1	C	18	GLY	4.9
1	A	59	PRO	4.9
2	D	214	LEU	4.9
1	A	134	SER	4.8
1	A	66	LYS	4.8
1	C	67	ASP	4.8
1	C	63	ILE	4.7
1	C	131	THR	4.7
1	C	70	LYS	4.7
1	A	136	ASN	4.6
1	C	222	GLN	4.5
1	C	28	GLU	4.5
1	A	19	PRO	4.4
2	D	361	HIS	4.4
1	C	138	GLU	4.4
1	A	21	VAL	4.3
1	A	290	THR	4.3
1	A	138	GLU	4.2
1	C	254	VAL	4.2
3	T	706	DA	4.2
1	A	251	SER	4.2
1	A	72	ARG	4.1
2	D	232	TYR	4.1
2	D	358	ARG	4.1
1	A	20	LYS	4.1
1	A	29	GLU	4.1
1	C	68	SER	4.0
1	A	293	ILE	4.0
1	C	139	THR	3.9
1	C	224	GLU	3.8
1	A	286	THR	3.8
1	A	69	THR	3.8
2	B	85	GLN	3.7
1	A	144	TYR	3.7
1	C	66	LYS	3.7
1	A	22	LYS	3.6
1	A	295	LEU	3.6
1	A	130	PHE	3.5
1	C	29	GLU	3.5
1	C	145	GLN	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	23	GLN	3.4
1	A	71	TRP	3.4
2	B	215	THR	3.4
2	D	212	TRP	3.4
1	C	52	PRO	3.4
1	A	219	LYS	3.3
2	D	215	THR	3.3
1	C	205	LEU	3.3
1	A	131	THR	3.3
1	A	70	LYS	3.3
3	T	725	DG	3.3
2	D	202	ILE	3.2
3	E	725	DG	3.2
1	C	19	PRO	3.1
2	D	359	GLY	3.1
1	C	130	PHE	3.1
1	C	193	LEU	3.1
3	E	706	DA	3.1
1	A	63	ILE	3.1
1	C	51	GLY	3.0
2	B	4	PRO	3.0
1	C	295	LEU	3.0
2	D	200	THR	3.0
1	C	128	THR	3.0
1	A	288	ALA	3.0
1	C	220	LYS	3.0
2	D	428	GLN	2.9
2	D	93	GLY	2.9
2	B	216	THR	2.9
1	C	141	GLY	2.9
1	C	76	ASP	2.9
1	C	22	LYS	2.8
4	P	803	DC	2.8
1	A	142	ILE	2.8
1	C	58	THR	2.8
2	D	89	GLU	2.8
3	T	724	DT	2.8
1	A	280	SER	2.8
2	D	88	TRP	2.8
1	C	109	LEU	2.8
1	A	283	LEU	2.7
1	C	50	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	88	TRP	2.7
1	A	141	GLY	2.7
1	A	221	HIS	2.7
1	C	46	LYS	2.7
2	B	86	ASP	2.7
1	C	55	PRO	2.7
1	C	30	LYS	2.7
2	D	357	MET	2.7
1	C	140	PRO	2.6
1	A	223	LYS	2.6
1	A	554	ALA	2.6
1	A	73	LYS	2.6
2	B	299	ALA	2.6
2	D	86	ASP	2.6
1	C	23	GLN	2.6
2	B	281	LYS	2.6
1	C	202	ILE	2.5
2	D	174	GLN	2.5
1	A	36	GLU	2.5
1	C	310	LEU	2.5
1	A	195	ILE	2.5
1	A	260	LEU	2.5
1	C	247	PRO	2.5
2	D	354	TYR	2.4
1	A	297	GLU	2.4
2	D	193	LEU	2.4
1	C	257	ILE	2.4
1	C	248	GLU	2.4
1	A	254	VAL	2.4
1	C	129	ALA	2.4
1	A	65	LYS	2.4
1	C	152	GLY	2.3
2	D	6	GLU	2.3
1	C	91	GLN	2.3
2	B	92	LEU	2.3
2	B	211	ARG	2.3
1	A	146	TYR	2.3
1	A	135	ILE	2.3
1	C	135	ILE	2.3
2	B	301	LEU	2.3
4	F	816	DG	2.2
1	C	448	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	189	VAL	2.2
1	A	15	GLY	2.2
1	C	346	PHE	2.2
1	C	142	ILE	2.2
1	C	286	THR	2.2
2	B	89	GLU	2.2
2	D	7	THR	2.2
1	A	147	ASN	2.2
1	C	13	LYS	2.1
1	A	82	LYS	2.1
1	C	75	VAL	2.1
1	C	21	VAL	2.1
1	A	292	VAL	2.1
1	C	146	TYR	2.1
1	C	283	LEU	2.1
1	A	285	GLY	2.1
2	D	66	LYS	2.1
2	B	422	LEU	2.1
1	A	284	ARG	2.1
1	A	56	TYR	2.0
2	B	94	ILE	2.0
1	A	282	LEU	2.0
1	A	217	PRO	2.0
1	C	20	LYS	2.0
2	B	212	TRP	2.0
1	C	189	VAL	2.0
1	C	285	GLY	2.0
2	B	231	GLY	2.0
1	C	14	PRO	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	MRG	P	817	26/27	0.86	0.27	-	87,107,136,143	0
4	ATM	P	822	22/23	0.91	0.18	-	103,116,127,129	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	MRG	F	817	26/27	0.81	0.31	-	97,121,156,164	0
4	ATM	F	822	22/23	0.84	0.23	-	107,118,133,136	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(\AA^2)	Q<0.9
5	NVP	C	901	20/20	0.95	0.26	0.76	67,73,78,88	0
5	NVP	A	901	20/20	0.94	0.23	0.12	63,69,75,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.