



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:03 AM UTC

PDB ID : 3KK3 / pdb_00003kk3
Title : HIV-1 reverse transcriptase-DNA complex with GS-9148 terminated primer
Authors : Lansdon, E.B.
Deposited on : 2009-11-04
Resolution : 2.90 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

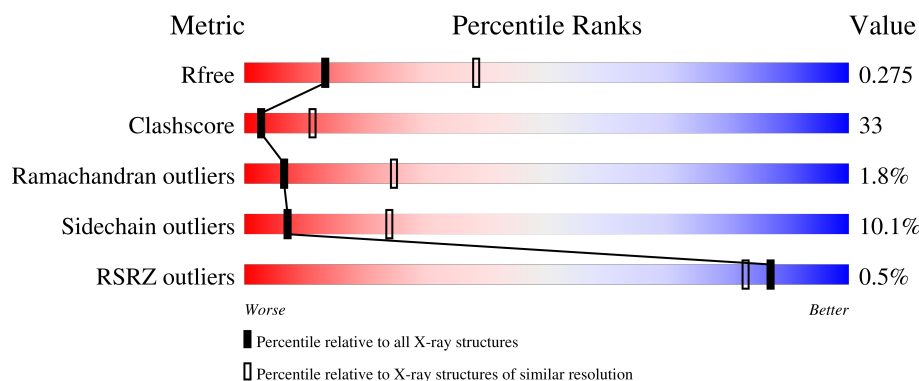
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	A	560	% 44% 47% 8% ..
2	B	452	35% 47% 8% 10%
3	P	21	5% 19% 62% 14%
4	T	27	. 30% 52% 15%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8752 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 Reverse transcriptase p66 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4514	2917	753	836	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	CYS	GLN	engineered mutation	UNP P04585
A	280	SER	CYS	engineered mutation	UNP P04585

- Molecule 2 is a protein called Reverse transcriptase p51 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	407	Total	C	N	O	S	0	0	0
			3361	2191	555	609	6			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	MET	-	expression tag	UNP P04585
B	-10	GLY	-	expression tag	UNP P04585
B	-9	SER	-	expression tag	UNP P04585
B	-8	SER	-	expression tag	UNP P04585
B	-7	HIS	-	expression tag	UNP P04585
B	-6	HIS	-	expression tag	UNP P04585
B	-5	HIS	-	expression tag	UNP P04585
B	-4	HIS	-	expression tag	UNP P04585
B	-3	HIS	-	expression tag	UNP P04585
B	-2	HIS	-	expression tag	UNP P04585
B	-1	SER	-	expression tag	UNP P04585
B	0	SER	-	expression tag	UNP P04585
B	280	SER	CYS	engineered mutation	UNP P04585

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	P	18	Total	C	F	N	O	P	0	0	0
			366	173	1	64	110	18			

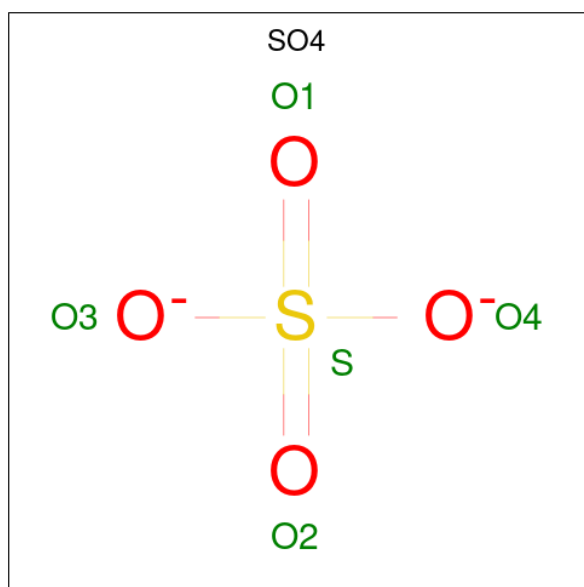
- Molecule 4 is a DNA chain called 5'-D(*AP*TP*GP*GP*TP*TP*GP*GP*CP*GP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*GP*TP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	T	23	Total	C	N	O	P	0	0	0
			478	224	94	137	23			

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	P	1	Total	O	S	0	0
			5	4	1		
6	T	1	Total	O	S	0	0
			5	4	1		

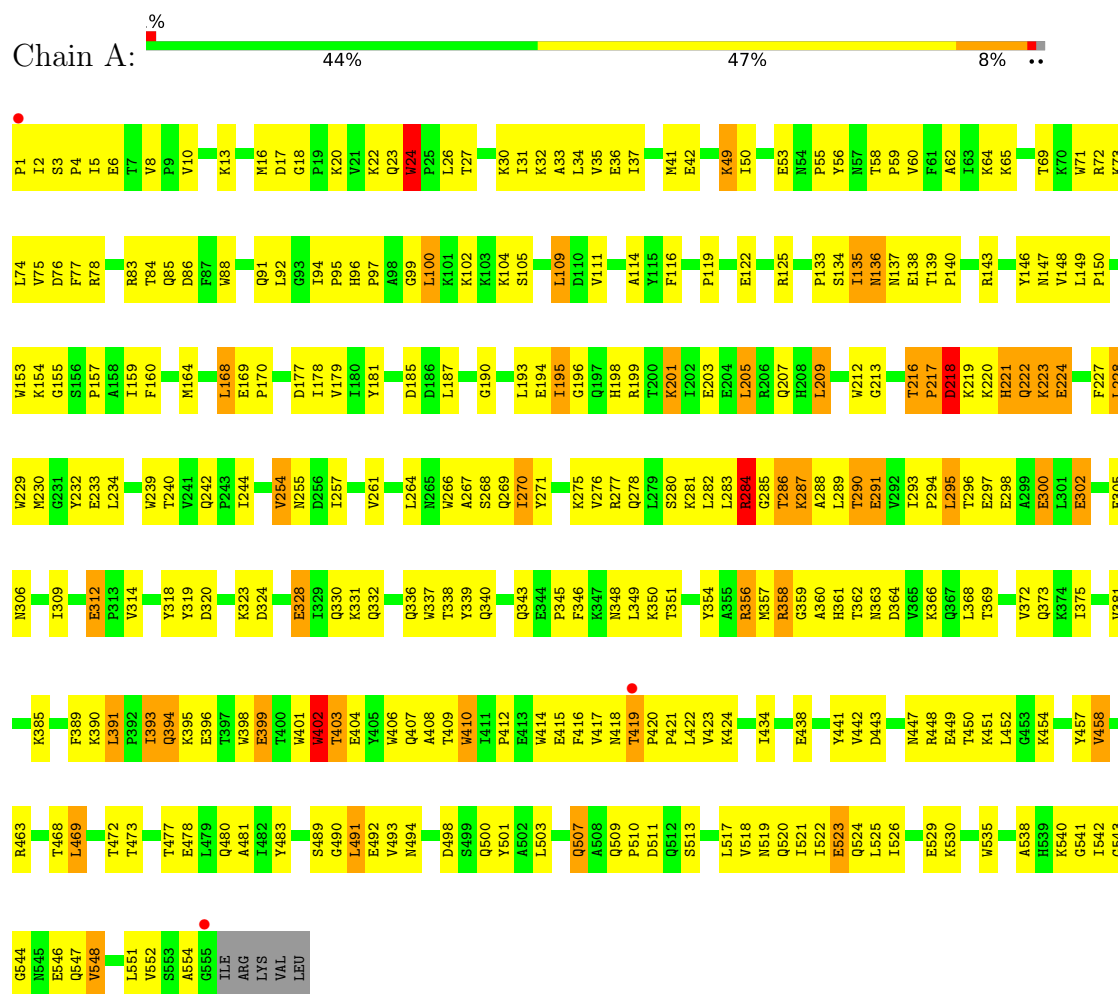
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	15	Total 15	O 15	0	0
7	B	4	Total 4	O 4	0	0
7	P	2	Total 2	O 2	0	0
7	T	1	Total 1	O 1	0	0

3 Residue-property plots

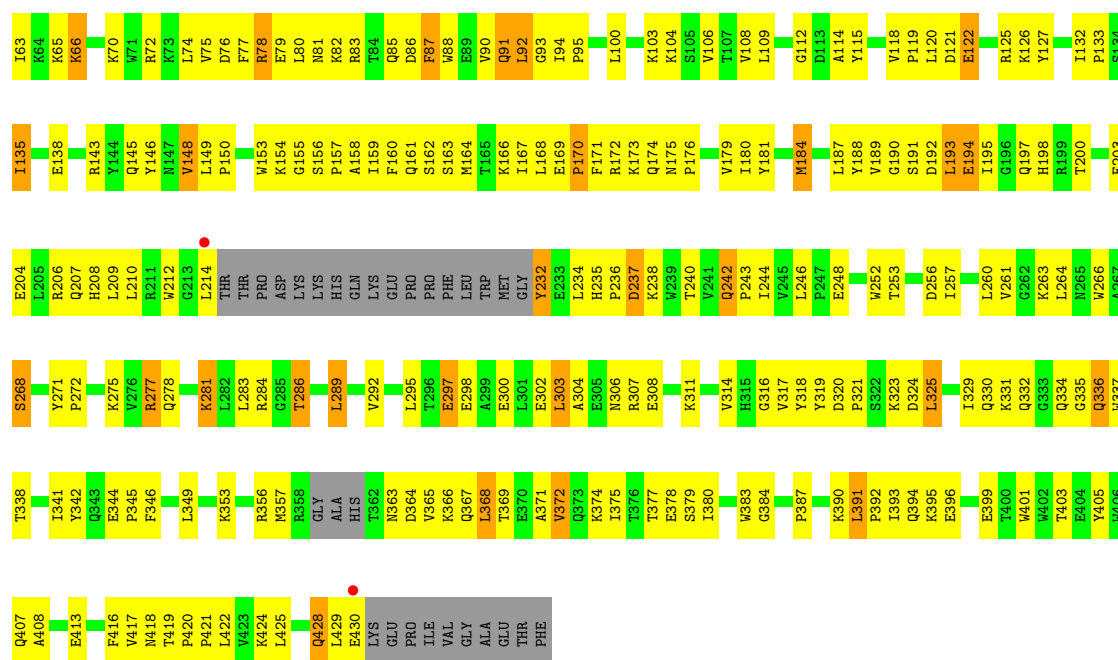
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Reverse transcriptase p66 subunit



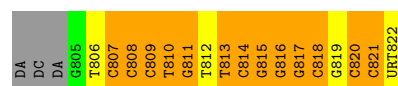
• Molecule 2: Reverse transcriptase p51 subunit





• Molecule 3: 5'-D(*AP*CP*AP*GP*TP*CP*CP*CP*TP*GP*TP*TP*CP*GP*GP*GP*CP*GP*CP*CP*(URT))-3'

Chain P: 5% 19% 62% 14%



• Molecule 4: 5'-D(*AP*TP*GP*GP*TP*TP*GP*GP*CP*GP*CP*CP*CP*GP*AP*AP*CP*AP*GP*GP*GP*AP*CP*TP*GP*TP*G))-3'

Chain T: 30% 52% 15%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	167.00Å 169.07Å 96.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.99 – 2.90 29.99 – 2.90	Depositor EDS
% Data completeness (in resolution range)	89.1 (29.99-2.90) 89.0 (29.99-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.50 (at 2.77Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.223 , 0.287 0.209 , 0.275	Depositor DCC
R_{free} test set	1343 reflections (4.27%)	wwPDB-VP
Wilson B-factor (Å ²)	68.9	Xtriage
Anisotropy	0.516	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8752	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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, 0.333 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: URT, MG, SO4

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	0.55	0/4631	0.91	13/6292 (0.2%)
2	B	0.46	0/3455	0.89	16/4694 (0.3%)
3	P	0.44	0/384	1.31	12/590 (2.0%)
4	T	0.40	0/537	1.21	17/828 (2.1%)
All	All	0.51	0/9007	0.95	58/12404 (0.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
3	P	0	1
All	All	0	4

There are no bond length outliers.

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	GLU	CA-C-N	-9.75	110.33	120.38
1	A	224	GLU	C-N-CA	-9.75	110.33	120.38
2	B	18	GLY	CA-C-N	8.96	129.31	119.90
2	B	18	GLY	C-N-CA	8.96	129.31	119.90
1	A	58	THR	CA-C-N	7.64	128.08	119.83
1	A	58	THR	C-N-CA	7.64	128.08	119.83
2	B	8	VAL	CA-C-N	7.62	127.38	119.76
2	B	8	VAL	C-N-CA	7.62	127.38	119.76
2	B	118	VAL	N-CA-C	7.51	116.29	107.73
4	T	723	DC	C2'-C3'-O3'	-7.18	100.73	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	816	DG	C2'-C3'-O3'	-6.97	101.05	111.50
3	P	821	DC	C2'-C3'-O3'	-6.91	101.13	111.50
4	T	714	DG	C2'-C3'-O3'	-6.82	101.27	111.50
4	T	720	DG	C2'-C3'-O3'	-6.66	101.51	111.50
3	P	809	DC	C2'-C3'-O3'	-6.60	101.59	111.50
4	T	721	DG	C2'-C3'-O3'	-6.39	101.92	111.50
2	B	132	ILE	CA-C-N	6.20	126.16	120.21
2	B	132	ILE	C-N-CA	6.20	126.16	120.21
1	A	24	TRP	CA-C-N	6.17	126.14	119.78
1	A	24	TRP	C-N-CA	6.17	126.14	119.78
3	P	818	DC	C2'-C3'-O3'	-6.15	102.27	111.50
4	T	713	DC	C2'-C3'-O3'	-6.03	102.45	111.50
1	A	360	ALA	N-CA-C	-5.91	105.91	113.23
3	P	813	DT	C2'-C3'-O3'	-5.87	102.70	111.50
3	P	815	DG	C2'-C3'-O3'	-5.85	102.73	111.50
4	T	711	DC	C2'-C3'-O3'	-5.80	102.79	111.50
2	B	54	ASN	CA-C-N	5.80	125.90	119.87
2	B	54	ASN	C-N-CA	5.80	125.90	119.87
4	T	708	DG	C2'-C3'-O3'	-5.75	102.87	111.50
3	P	817	DG	C2'-C3'-O3'	-5.71	102.93	111.50
4	T	724	DT	C2'-C3'-O3'	-5.71	102.94	111.50
3	P	810	DT	C2'-C3'-O3'	-5.70	102.95	111.50
1	A	242	GLN	CA-C-N	5.62	125.64	119.90
1	A	242	GLN	C-N-CA	5.62	125.64	119.90
1	A	410	TRP	N-CA-C	5.60	117.58	109.07
4	T	705	DT	C2'-C3'-O3'	-5.58	103.13	111.50
2	B	235	HIS	CA-C-N	5.51	126.73	119.84
2	B	235	HIS	C-N-CA	5.51	126.73	119.84
4	T	719	DG	C2'-C3'-O3'	-5.47	103.30	111.50
4	T	716	DA	C2'-C3'-O3'	-5.46	103.30	111.50
4	T	715	DA	C2'-C3'-O3'	-5.38	103.44	111.50
2	B	391	LEU	CA-C-N	5.35	128.50	120.96
2	B	391	LEU	C-N-CA	5.35	128.50	120.96
2	B	7	THR	N-CA-C	5.32	117.49	109.41
2	B	401	TRP	N-CA-C	5.29	121.48	114.12
4	T	709	DC	C2'-C3'-O3'	-5.28	103.58	111.50
3	P	808	DC	C2'-C3'-O3'	-5.23	103.65	111.50
4	T	717	DC	C2'-C3'-O3'	-5.23	103.66	111.50
3	P	814	DC	C2'-C3'-O3'	-5.21	103.69	111.50
2	B	184	MET	CB-CA-C	-5.18	109.62	115.79
4	T	706	DT	C2'-C3'-O3'	-5.17	103.75	111.50
1	A	99	GLY	N-CA-C	-5.16	107.61	115.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	222	GLN	N-CA-C	-5.13	102.16	110.32
3	P	811	DG	C2'-C3'-O3'	-5.12	103.83	111.50
1	A	402	TRP	N-CA-C	-5.09	106.76	112.87
3	P	807	DC	C2'-C3'-O3'	-5.06	103.91	111.50
4	T	722	DA	C2'-C3'-O3'	-5.03	103.96	111.50
4	T	714	DG	N9-C1'-C2'	5.01	121.02	113.50

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	218	ASP	Peptide
1	A	286	THR	Peptide
1	A	287	LYS	Peptide
3	P	820	DC	Sidechain

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	4514	0	4563	310	0
2	B	3361	0	3393	232	1
3	P	366	0	200	26	0
4	T	478	0	257	32	0
5	A	1	0	0	0	0
6	P	5	0	0	0	0
6	T	5	0	0	0	0
7	A	15	0	0	1	0
7	B	4	0	0	1	0
7	P	2	0	0	0	0
7	T	1	0	0	0	0
All	All	8752	0	8413	569	1

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 33.

All (569) 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:336:GLN:HE22	2:B:353:LYS:HD2	1.26	1.00
1:A:125:ARG:HG2	1:A:146:TYR:O	1.63	0.98
2:B:5:ILE:HG23	2:B:6:GLU:H	1.30	0.96
1:A:195:ILE:HD13	1:A:195:ILE:H	1.29	0.96
1:A:255:ASN:HD22	1:A:289:LEU:HD22	1.28	0.96
1:A:37:ILE:HG22	1:A:41:MET:HE2	1.49	0.93
3:P:820:DC:H2''	3:P:821:DC:H5'	1.48	0.92
2:B:169:GLU:HB3	2:B:170:PRO:HD3	1.53	0.90
2:B:35:VAL:O	2:B:39:THR:HG23	1.73	0.89
1:A:20:LYS:HE2	1:A:55:PRO:HB2	1.56	0.87
2:B:277:ARG:NH1	2:B:281:LYS:HD3	1.91	0.85
1:A:199:ARG:NH2	1:A:223:LYS:HB2	1.90	0.85
2:B:180:ILE:HG23	2:B:189:VAL:HG22	1.59	0.85
1:A:136:ASN:ND2	1:A:138:GLU:HB3	1.91	0.85
2:B:429:LEU:HD23	2:B:430:GLU:H	1.40	0.84
4:T:713:DC:H2'	4:T:714:DG:C8	2.13	0.83
1:A:518:VAL:O	1:A:522:ILE:HG13	1.81	0.81
1:A:143:ARG:HG3	1:A:143:ARG:HH11	1.45	0.81
2:B:86:ASP:HA	2:B:88:TRP:CZ3	2.15	0.81
1:A:296:THR:HG22	1:A:298:GLU:H	1.45	0.80
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.62	0.80
1:A:297:GLU:HA	1:A:300:GLU:HB2	1.65	0.79
1:A:53:GLU:H	1:A:53:GLU:CD	1.91	0.79
2:B:122:GLU:HA	2:B:125:ARG:HH11	1.48	0.78
1:A:34:LEU:HD21	1:A:62:ALA:HB2	1.66	0.78
1:A:224:GLU:O	1:A:224:GLU:HG2	1.81	0.77
1:A:164:MET:HE3	1:A:168:LEU:HD21	1.66	0.77
2:B:420:PRO:HB2	2:B:422:LEU:HD23	1.67	0.77
1:A:364:ASP:HB3	1:A:423:VAL:HG13	1.67	0.76
2:B:103:LYS:HD2	2:B:191:SER:HA	1.68	0.75
3:P:813:DT:H2'	3:P:814:DC:C6	2.22	0.75
1:A:293:ILE:HD12	1:A:294:PRO:HD2	1.68	0.75
1:A:30:LYS:HG2	1:A:71:TRP:CZ3	2.22	0.75
1:A:109:LEU:HD12	1:A:187:LEU:HB2	1.68	0.75
2:B:345:PRO:O	2:B:346:PHE:HB2	1.86	0.75
1:A:295:LEU:HG	1:A:300:GLU:OE2	1.86	0.75
2:B:56:TYR:O	2:B:143:ARG:NH2	2.19	0.74
1:A:356:ARG:HD2	1:A:358:ARG:HG3	1.68	0.74
2:B:115:TYR:OH	2:B:157:PRO:HB3	1.88	0.74
1:A:31:ILE:O	1:A:35:VAL:HG23	1.88	0.73
1:A:275:LYS:H	1:A:306:ASN:HD21	1.34	0.73
2:B:65:LYS:O	2:B:66:LYS:HD2	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:86:ASP:HA	2:B:88:TRP:HZ3	1.52	0.73
2:B:126:LYS:HA	2:B:145:GLN:OE1	1.87	0.73
1:A:138:GLU:HG2	1:A:139:THR:N	2.03	0.73
1:A:417:VAL:HG12	1:A:419:THR:HG22	1.70	0.72
2:B:10:VAL:HG12	2:B:11:LYS:H	1.54	0.72
1:A:26:LEU:HD12	1:A:133:PRO:HG3	1.70	0.72
1:A:358:ARG:HB3	1:A:358:ARG:HH11	1.53	0.72
1:A:363:ASN:HA	1:A:511:ASP:OD2	1.90	0.70
1:A:520:GLN:O	1:A:523:GLU:HG3	1.92	0.70
4:T:703:DG:H2''	4:T:704:DG:N2	2.07	0.70
1:A:2:ILE:CG2	1:A:119:PRO:HG3	2.23	0.69
2:B:79:GLU:HG3	2:B:83:ARG:HH11	1.57	0.69
1:A:8:VAL:HG21	1:A:159:ILE:HG23	1.75	0.68
1:A:97:PRO:O	1:A:100:LEU:HB2	1.94	0.68
1:A:278:GLN:HB2	1:A:302:GLU:CD	2.18	0.68
2:B:200:THR:O	2:B:204:GLU:HG3	1.93	0.68
2:B:275:LYS:HE3	2:B:277:ARG:HB2	1.75	0.68
2:B:163:SER:O	2:B:167:ILE:HG13	1.95	0.67
1:A:277:ARG:HB3	1:A:336:GLN:NE2	2.09	0.67
1:A:418:ASN:O	1:A:420:PRO:HD3	1.94	0.67
1:A:492:GLU:OE1	1:A:530:LYS:HD2	1.94	0.67
2:B:331:LYS:HB2	2:B:337:TRP:CZ3	2.29	0.67
2:B:429:LEU:HD23	2:B:430:GLU:N	2.09	0.67
1:A:65:LYS:HG2	1:A:72:ARG:HE	1.58	0.67
2:B:82:LYS:HE2	2:B:413:GLU:OE2	1.94	0.67
4:T:704:DG:H2''	4:T:705:DT:OP1	1.94	0.67
4:T:713:DC:H2''	4:T:714:DG:O4'	1.94	0.67
1:A:255:ASN:ND2	1:A:289:LEU:HD22	2.05	0.67
2:B:5:ILE:HG23	2:B:6:GLU:N	2.08	0.67
2:B:275:LYS:HE3	2:B:277:ARG:CB	2.25	0.66
2:B:336:GLN:NE2	2:B:353:LYS:HD2	2.05	0.66
1:A:403:THR:HG22	1:A:404:GLU:N	2.11	0.66
2:B:317:VAL:HG12	2:B:349:LEU:HD23	1.79	0.65
2:B:31:ILE:HD12	2:B:135:ILE:HG22	1.79	0.65
2:B:278:GLN:HB2	2:B:302:GLU:OE1	1.96	0.65
1:A:122:GLU:HG2	1:A:125:ARG:NH2	2.11	0.65
1:A:366:LYS:O	1:A:369:THR:HB	1.96	0.65
2:B:13:LYS:HZ1	2:B:88:TRP:HH2	1.45	0.65
1:A:199:ARG:HH21	1:A:223:LYS:HB2	1.59	0.65
2:B:104:LYS:HB3	2:B:192:ASP:HA	1.78	0.65
1:A:390:LYS:HB3	1:A:417:VAL:HG21	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:707:DG:H2''	4:T:708:DG:H5'	1.79	0.65
1:A:478:GLU:HA	1:A:478:GLU:OE1	1.96	0.65
2:B:180:ILE:HG12	2:B:189:VAL:HG13	1.79	0.65
2:B:206:ARG:O	2:B:210:LEU:HD13	1.97	0.65
1:A:60:VAL:HG21	1:A:73:LYS:HE2	1.78	0.65
2:B:425:LEU:HD12	2:B:428:GLN:HG3	1.79	0.65
1:A:222:GLN:O	1:A:223:LYS:C	2.38	0.64
2:B:109:LEU:HB3	2:B:187:LEU:HB3	1.79	0.64
4:T:708:DG:H2'	4:T:709:DC:C6	2.33	0.64
1:A:22:LYS:O	1:A:22:LYS:HG2	1.98	0.64
1:A:395:LYS:HD3	1:A:414:TRP:CZ2	2.32	0.64
1:A:164:MET:HE3	1:A:168:LEU:CD2	2.29	0.63
1:A:195:ILE:HD13	1:A:195:ILE:N	2.10	0.63
2:B:122:GLU:HA	2:B:125:ARG:NH1	2.13	0.63
2:B:26:LEU:HD12	2:B:133:PRO:CG	2.28	0.63
4:T:718:DA:H2''	4:T:719:DG:OP1	1.97	0.63
1:A:277:ARG:HB3	1:A:336:GLN:CD	2.23	0.63
2:B:33:ALA:O	2:B:37:ILE:HG13	1.98	0.63
2:B:88:TRP:O	2:B:92:LEU:HD12	1.97	0.63
1:A:27:THR:OG1	1:A:30:LYS:HG3	1.98	0.63
1:A:30:LYS:HE2	4:T:703:DG:H22	1.64	0.63
1:A:33:ALA:O	1:A:37:ILE:HG13	1.98	0.63
1:A:393:ILE:HG13	1:A:423:VAL:HB	1.80	0.63
2:B:170:PRO:HB2	2:B:208:HIS:NE2	2.14	0.63
2:B:337:TRP:HE1	2:B:367:GLN:HG2	1.64	0.63
1:A:500:GLN:HE21	2:B:420:PRO:CB	2.12	0.62
2:B:172:ARG:NH2	2:B:180:ILE:HB	2.14	0.62
1:A:143:ARG:HG3	1:A:143:ARG:NH1	2.14	0.62
1:A:37:ILE:CG2	1:A:41:MET:HE2	2.28	0.62
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.82	0.62
2:B:81:ASN:O	2:B:154:LYS:HE3	2.00	0.62
2:B:277:ARG:HH11	2:B:281:LYS:HD3	1.65	0.62
1:A:94:ILE:CD1	4:T:709:DC:H1'	2.29	0.61
1:A:390:LYS:HB3	1:A:417:VAL:CG2	2.30	0.61
1:A:94:ILE:HD13	4:T:709:DC:H1'	1.82	0.61
1:A:135:ILE:HG23	1:A:136:ASN:CG	2.25	0.61
2:B:109:LEU:HD23	2:B:187:LEU:HD23	1.82	0.61
4:T:714:DG:H2'	4:T:715:DA:C8	2.36	0.61
1:A:1:PRO:HG2	1:A:213:GLY:HA2	1.83	0.61
2:B:422:LEU:N	2:B:422:LEU:HD22	2.16	0.61
2:B:368:LEU:O	2:B:368:LEU:HD22	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:197:GLN:HA	2:B:200:THR:HG22	1.83	0.60
1:A:1:PRO:O	1:A:2:ILE:HD13	2.01	0.60
2:B:79:GLU:HG3	2:B:83:ARG:NH1	2.16	0.60
1:A:285:GLY:O	1:A:287:LYS:HB3	2.01	0.60
1:A:104:LYS:HD2	1:A:193:LEU:O	2.01	0.60
1:A:361:HIS:ND1	1:A:513:SER:HB2	2.16	0.60
1:A:399:GLU:HA	1:A:402:TRP:HD1	1.67	0.60
1:A:469:LEU:HB2	1:A:472:THR:HG21	1.83	0.60
2:B:203:GLU:OE1	2:B:206:ARG:HD3	2.02	0.60
2:B:244:ILE:N	2:B:244:ILE:HD12	2.16	0.60
1:A:23:GLN:NE2	1:A:59:PRO:HA	2.17	0.60
1:A:442:VAL:HB	1:A:481:ALA:HB1	1.83	0.60
2:B:316:GLY:HA2	2:B:318:TYR:CE2	2.37	0.60
1:A:458:VAL:HG13	2:B:286:THR:HG21	1.83	0.60
2:B:31:ILE:CD1	2:B:135:ILE:HG22	2.32	0.60
2:B:100:LEU:O	2:B:100:LEU:HD12	2.02	0.60
2:B:344:GLU:OE1	2:B:344:GLU:HA	2.01	0.60
4:T:720:DG:H2'	4:T:721:DG:C8	2.37	0.60
1:A:254:VAL:O	1:A:257:ILE:HB	2.02	0.59
1:A:169:GLU:HB3	1:A:170:PRO:HD3	1.83	0.59
2:B:90:VAL:HG12	2:B:91:GLN:HE22	1.67	0.59
2:B:87:PHE:CZ	2:B:155:GLY:HA2	2.37	0.59
1:A:16:MET:HB3	1:A:83:ARG:NH1	2.17	0.59
1:A:23:GLN:HE22	1:A:59:PRO:HA	1.67	0.59
1:A:199:ARG:NH2	1:A:223:LYS:CB	2.62	0.59
2:B:257:ILE:O	2:B:261:VAL:HG23	2.03	0.59
1:A:320:ASP:OD2	1:A:323:LYS:HG3	2.03	0.59
1:A:543:GLY:HA3	2:B:283:LEU:O	2.03	0.58
2:B:104:LYS:O	2:B:236:PRO:HD2	2.04	0.58
2:B:421:PRO:O	2:B:424:LYS:HB3	2.03	0.58
3:P:817:DG:H2'	3:P:818:DC:C6	2.37	0.58
3:P:820:DC:H2''	3:P:821:DC:C5'	2.27	0.58
2:B:335:GLY:HA2	2:B:367:GLN:NE2	2.18	0.58
1:A:64:LYS:HE3	1:A:71:TRP:CZ2	2.39	0.58
1:A:135:ILE:HG23	1:A:136:ASN:OD1	2.02	0.58
2:B:5:ILE:HG12	2:B:6:GLU:HG3	1.84	0.58
1:A:195:ILE:H	1:A:195:ILE:CD1	2.09	0.58
2:B:191:SER:H	2:B:198:HIS:HE1	1.52	0.58
3:P:818:DC:H2''	3:P:819:DG:H5'	1.85	0.58
1:A:60:VAL:O	1:A:60:VAL:HG13	2.03	0.58
1:A:295:LEU:HD12	1:A:296:THR:O	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:THR:O	1:A:69:THR:HG22	2.03	0.57
1:A:522:ILE:O	1:A:526:ILE:HG13	2.04	0.57
2:B:160:PHE:O	2:B:164:MET:HB2	2.04	0.57
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.40	0.57
1:A:469:LEU:N	1:A:469:LEU:HD22	2.20	0.57
2:B:90:VAL:HG12	2:B:91:GLN:NE2	2.20	0.57
2:B:191:SER:HB2	2:B:193:LEU:HD13	1.85	0.57
1:A:420:PRO:HA	1:A:421:PRO:C	2.29	0.57
2:B:277:ARG:O	2:B:277:ARG:HD3	2.05	0.57
1:A:394:GLN:HB2	1:A:396:GLU:OE1	2.04	0.57
2:B:108:VAL:HG13	2:B:188:TYR:CE2	2.40	0.57
1:A:136:ASN:HD21	1:A:138:GLU:HB3	1.70	0.57
2:B:13:LYS:HB2	2:B:16:MET:HG3	1.85	0.57
1:A:139:THR:HG22	1:A:140:PRO:O	2.05	0.57
2:B:112:GLY:C	2:B:114:ALA:H	2.10	0.57
1:A:398:TRP:CH2	1:A:409:THR:HG23	2.39	0.56
2:B:368:LEU:O	2:B:372:VAL:HG13	2.05	0.56
2:B:390:LYS:HB3	2:B:417:VAL:HG21	1.86	0.56
1:A:473:THR:O	1:A:477:THR:HG23	2.04	0.56
2:B:92:LEU:HD22	2:B:94:ILE:HG13	1.86	0.56
1:A:85:GLN:HG3	1:A:86:ASP:O	2.05	0.56
1:A:94:ILE:HD12	1:A:94:ILE:N	2.20	0.56
1:A:114:ALA:HB1	1:A:160:PHE:CZ	2.40	0.56
1:A:328:GLU:HG2	1:A:390:LYS:HB2	1.86	0.56
2:B:209:LEU:O	2:B:212:TRP:N	2.38	0.56
1:A:332:GLN:HG3	1:A:336:GLN:O	2.05	0.56
2:B:103:LYS:HE2	2:B:179:VAL:HG12	1.87	0.56
2:B:169:GLU:HB3	2:B:170:PRO:CD	2.30	0.56
1:A:30:LYS:CE	4:T:703:DG:N2	2.68	0.56
1:A:233:GLU:HB2	1:A:240:THR:HG23	1.87	0.56
3:P:810:DT:H2'	3:P:811:DG:C8	2.41	0.56
2:B:356:ARG:HD2	2:B:367:GLN:HG3	1.87	0.56
1:A:136:ASN:HD22	1:A:138:GLU:HB3	1.71	0.55
1:A:216:THR:O	1:A:217:PRO:O	2.24	0.55
1:A:389:PHE:HB3	1:A:391:LEU:CD1	2.36	0.55
1:A:547:GLN:O	1:A:551:LEU:HD13	2.06	0.55
1:A:406:TRP:HA	2:B:331:LYS:HD2	1.88	0.55
2:B:88:TRP:HB3	2:B:154:LYS:HD2	1.89	0.55
2:B:194:GLU:HG2	2:B:195:ILE:N	2.22	0.55
3:P:813:DT:H2'	3:P:814:DC:H6	1.67	0.55
1:A:205:LEU:HD22	1:A:209:LEU:HD22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:LYS:O	1:A:223:LYS:HG2	2.06	0.55
2:B:7:THR:HB	2:B:119:PRO:HB2	1.89	0.54
1:A:230:MET:HA	3:P:821:DC:O3'	2.07	0.54
2:B:308:GLU:O	2:B:311:LYS:HG3	2.08	0.54
1:A:30:LYS:HE2	4:T:703:DG:N2	2.22	0.54
2:B:10:VAL:HG12	2:B:11:LYS:N	2.21	0.54
2:B:11:LYS:HB3	2:B:85:GLN:OE1	2.08	0.54
4:T:715:DA:H2'	4:T:716:DA:C8	2.42	0.54
2:B:379:SER:HA	2:B:383:TRP:CE3	2.42	0.54
2:B:422:LEU:HD22	2:B:422:LEU:H	1.71	0.54
3:P:813:DT:H2''	3:P:814:DC:H5'	1.90	0.54
1:A:97:PRO:HA	1:A:100:LEU:HD22	1.90	0.54
2:B:159:ILE:C	2:B:161:GLN:H	2.15	0.54
1:A:257:ILE:HD12	1:A:293:ILE:HG21	1.89	0.53
2:B:345:PRO:O	2:B:346:PHE:CB	2.55	0.53
1:A:31:ILE:HD13	1:A:135:ILE:HA	1.91	0.53
1:A:510:PRO:HB2	1:A:522:ILE:HD11	1.89	0.53
2:B:115:TYR:CD2	2:B:156:SER:HB3	2.43	0.53
2:B:120:LEU:O	2:B:125:ARG:NH1	2.42	0.53
4:T:721:DG:H2''	4:T:722:DA:O5'	2.08	0.53
1:A:32:LYS:O	1:A:36:GLU:HG3	2.09	0.53
2:B:28:GLU:OE1	2:B:32:LYS:HE2	2.08	0.53
1:A:199:ARG:HH21	1:A:223:LYS:CB	2.20	0.53
2:B:87:PHE:HB2	2:B:91:GLN:HB2	1.89	0.53
1:A:73:LYS:HG3	1:A:74:LEU:N	2.24	0.53
2:B:61:PHE:HD1	2:B:61:PHE:H	1.56	0.53
1:A:178:ILE:HD11	1:A:201:LYS:HG2	1.90	0.53
1:A:320:ASP:HB3	1:A:323:LYS:HZ2	1.75	0.52
2:B:24:TRP:HB2	2:B:25:PRO:HD2	1.90	0.52
1:A:448:ARG:HH12	4:T:723:DC:H1'	1.74	0.52
1:A:135:ILE:HD13	1:A:135:ILE:O	2.10	0.52
2:B:395:LYS:HE2	2:B:399:GLU:OE1	2.10	0.52
2:B:369:THR:HG21	2:B:405:TYR:HB2	1.90	0.52
1:A:134:SER:O	1:A:136:ASN:N	2.42	0.52
1:A:270:ILE:HG23	1:A:314:VAL:HG21	1.90	0.52
1:A:498:ASP:HB2	1:A:538:ALA:HB2	1.91	0.52
3:P:818:DC:H2'	3:P:819:DG:H8	1.73	0.52
1:A:296:THR:HG22	1:A:298:GLU:N	2.20	0.52
1:A:233:GLU:HB2	1:A:240:THR:CG2	2.39	0.51
2:B:12:LEU:O	2:B:13:LYS:C	2.53	0.51
2:B:194:GLU:HG2	2:B:195:ILE:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:341:ILE:N	2:B:341:ILE:HD12	2.25	0.51
1:A:393:ILE:O	1:A:414:TRP:CH2	2.62	0.51
1:A:434:ILE:HD12	1:A:493:VAL:O	2.11	0.51
2:B:366:LYS:O	2:B:369:THR:HG22	2.10	0.51
1:A:222:GLN:O	1:A:224:GLU:HB2	2.11	0.51
1:A:283:LEU:O	1:A:285:GLY:N	2.40	0.51
2:B:191:SER:H	2:B:198:HIS:CE1	2.28	0.51
3:P:817:DG:H2'	3:P:818:DC:H6	1.75	0.51
3:P:817:DG:H2''	3:P:818:DC:H5'	1.93	0.51
1:A:255:ASN:HD22	1:A:289:LEU:CD2	2.14	0.51
1:A:277:ARG:H	1:A:336:GLN:NE2	2.09	0.51
2:B:371:ALA:O	2:B:375:ILE:HG13	2.11	0.51
1:A:410:TRP:HB2	2:B:365:VAL:HG23	1.91	0.51
2:B:79:GLU:HB2	7:B:442:HOH:O	2.11	0.51
2:B:191:SER:CB	2:B:193:LEU:HD13	2.40	0.51
2:B:236:PRO:C	2:B:238:LYS:H	2.18	0.51
3:P:818:DC:H2'	3:P:819:DG:C8	2.46	0.51
3:P:819:DG:H2'	3:P:820:DC:C6	2.46	0.51
1:A:75:VAL:HB	1:A:77:PHE:CE2	2.46	0.50
1:A:454:LYS:CE	1:A:554:ALA:HB3	2.41	0.50
1:A:454:LYS:NZ	1:A:554:ALA:HB3	2.26	0.50
1:A:350:LYS:NZ	1:A:351:THR:O	2.45	0.50
2:B:66:LYS:HG2	2:B:232:TYR:CE1	2.47	0.50
2:B:319:TYR:CZ	2:B:321:PRO:HA	2.47	0.50
1:A:407:GLN:OE1	2:B:419:THR:HG22	2.12	0.50
2:B:380:ILE:O	2:B:384:GLY:HA2	2.12	0.50
1:A:222:GLN:O	1:A:224:GLU:CB	2.60	0.50
1:A:403:THR:HG22	1:A:404:GLU:HG3	1.92	0.50
2:B:334:GLN:NE2	2:B:334:GLN:HA	2.27	0.50
1:A:503:LEU:O	1:A:507:GLN:HB2	2.12	0.50
1:A:228:LEU:N	1:A:228:LEU:HD23	2.27	0.49
1:A:271:TYR:CE1	1:A:314:VAL:HG22	2.47	0.49
1:A:491:LEU:HG	1:A:529:GLU:HG3	1.94	0.49
1:A:179:VAL:HG12	1:A:190:GLY:O	2.12	0.49
2:B:332:GLN:HG3	2:B:338:THR:HG23	1.95	0.49
1:A:3:SER:C	1:A:5:ILE:H	2.20	0.49
1:A:278:GLN:HB2	1:A:302:GLU:OE1	2.13	0.49
2:B:252:TRP:O	2:B:292:VAL:HG13	2.12	0.49
2:B:253:THR:H	2:B:256:ASP:HB2	1.78	0.49
2:B:275:LYS:HE3	2:B:277:ARG:HB3	1.94	0.49
4:T:706:DT:H2'	4:T:707:DG:C8	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:ILE:HG12	1:A:6:GLU:N	2.27	0.49
1:A:297:GLU:CA	1:A:300:GLU:HB2	2.40	0.49
1:A:438:GLU:OE2	1:A:463:ARG:NH1	2.45	0.49
2:B:103:LYS:O	2:B:236:PRO:HG2	2.11	0.49
1:A:458:VAL:HG12	1:A:548:VAL:HG22	1.92	0.49
2:B:390:LYS:HB3	2:B:417:VAL:CG2	2.43	0.49
1:A:134:SER:CB	1:A:139:THR:HB	2.43	0.49
1:A:281:LYS:C	1:A:283:LEU:N	2.70	0.49
1:A:17:ASP:O	1:A:83:ARG:HD3	2.13	0.49
1:A:229:TRP:HB3	1:A:234:LEU:HD11	1.95	0.49
1:A:401:TRP:C	1:A:403:THR:N	2.70	0.49
2:B:27:THR:OG1	2:B:30:LYS:HG3	2.13	0.49
2:B:112:GLY:C	2:B:114:ALA:N	2.71	0.49
1:A:220:LYS:HE3	1:A:222:GLN:HG3	1.95	0.48
1:A:389:PHE:HB3	1:A:391:LEU:HD13	1.95	0.48
1:A:509:GLN:N	1:A:510:PRO:HD3	2.28	0.48
1:A:8:VAL:O	1:A:10:VAL:HG23	2.13	0.48
1:A:354:TYR:OH	1:A:356:ARG:HB3	2.13	0.48
1:A:153:TRP:O	1:A:155:GLY:N	2.46	0.48
1:A:410:TRP:CZ2	2:B:363:ASN:ND2	2.81	0.48
2:B:79:GLU:CG	2:B:83:ARG:NH1	2.76	0.48
1:A:50:ILE:HG13	1:A:143:ARG:HB3	1.94	0.48
1:A:195:ILE:HG12	1:A:196:GLY:H	1.79	0.48
1:A:348:ASN:O	1:A:349:LEU:C	2.57	0.48
1:A:401:TRP:C	1:A:403:THR:H	2.21	0.48
1:A:138:GLU:CG	1:A:139:THR:N	2.75	0.48
1:A:410:TRP:CZ2	1:A:412:PRO:HA	2.49	0.48
1:A:102:LYS:HB2	1:A:102:LYS:HE3	1.55	0.48
1:A:227:PHE:C	1:A:228:LEU:HD23	2.38	0.48
4:T:721:DG:C2'	4:T:722:DA:O5'	2.62	0.48
1:A:261:VAL:HG13	1:A:276:VAL:HG21	1.95	0.48
1:A:483:TYR:HB2	1:A:521:ILE:HG12	1.95	0.48
2:B:164:MET:HA	2:B:167:ILE:HD12	1.96	0.48
1:A:287:LYS:HE3	1:A:289:LEU:CD2	2.43	0.48
1:A:359:GLY:HA2	3:P:811:DG:OP2	2.14	0.48
1:A:393:ILE:O	1:A:414:TRP:CZ3	2.67	0.48
1:A:280:SER:HB2	4:T:712:DC:H4'	1.96	0.48
1:A:320:ASP:OD2	1:A:323:LYS:HE3	2.14	0.48
1:A:372:VAL:HG13	1:A:389:PHE:CZ	2.49	0.48
1:A:418:ASN:O	1:A:420:PRO:CD	2.62	0.48
1:A:517:LEU:HA	1:A:520:GLN:OE1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:66:LYS:HE2	2:B:232:TYR:HD1	1.78	0.47
3:P:813:DT:H2''	3:P:814:DC:C5'	2.43	0.47
1:A:203:GLU:O	1:A:207:GLN:HG3	2.13	0.47
2:B:76:ASP:OD1	2:B:78:ARG:HB2	2.14	0.47
1:A:218:ASP:OD1	1:A:218:ASP:N	2.46	0.47
1:A:324:ASP:O	1:A:343:GLN:HG2	2.14	0.47
2:B:325:LEU:HB3	2:B:387:PRO:HA	1.95	0.47
2:B:332:GLN:HB2	2:B:336:GLN:O	2.14	0.47
1:A:447:ASN:O	1:A:449:GLU:N	2.46	0.47
3:P:809:DC:H2''	3:P:810:DT:O5'	2.15	0.47
1:A:26:LEU:HD12	1:A:133:PRO:CG	2.40	0.47
2:B:166:LYS:O	2:B:168:LEU:N	2.46	0.47
1:A:266:TRP:O	1:A:269:GLN:HG2	2.14	0.47
1:A:410:TRP:CB	2:B:365:VAL:HG23	2.45	0.47
2:B:106:VAL:HG23	2:B:236:PRO:HG3	1.97	0.47
1:A:328:GLU:O	1:A:339:TYR:HA	2.15	0.47
2:B:106:VAL:HA	2:B:190:GLY:HA3	1.96	0.47
2:B:207:GLN:C	2:B:209:LEU:N	2.72	0.47
3:P:816:DG:H2''	3:P:817:DG:O5'	2.14	0.47
1:A:17:ASP:CG	1:A:18:GLY:H	2.23	0.47
2:B:31:ILE:O	2:B:35:VAL:HG13	2.15	0.47
4:T:704:DG:C2'	4:T:705:DT:OP1	2.63	0.47
1:A:17:ASP:OD1	1:A:56:TYR:OH	2.32	0.46
1:A:27:THR:H	1:A:30:LYS:HD2	1.80	0.46
1:A:125:ARG:HD3	1:A:147:ASN:HA	1.97	0.46
2:B:87:PHE:HZ	2:B:155:GLY:HA2	1.80	0.46
2:B:316:GLY:HA2	2:B:318:TYR:HE2	1.77	0.46
4:T:703:DG:H2''	4:T:704:DG:H22	1.80	0.46
1:A:289:LEU:CD1	3:P:817:DG:H4'	2.45	0.46
2:B:23:GLN:NE2	2:B:24:TRP:O	2.48	0.46
2:B:125:ARG:HG2	2:B:146:TYR:O	2.15	0.46
1:A:270:ILE:HG23	1:A:314:VAL:CG2	2.46	0.46
1:A:320:ASP:CB	1:A:323:LYS:NZ	2.78	0.46
1:A:501:TYR:CE1	3:P:809:DC:H4'	2.51	0.46
2:B:16:MET:HE3	2:B:83:ARG:HG2	1.97	0.46
2:B:337:TRP:HE1	2:B:367:GLN:HE21	1.63	0.46
1:A:254:VAL:CG1	1:A:289:LEU:HA	2.46	0.46
2:B:393:ILE:HG12	2:B:394:GLN:N	2.30	0.46
1:A:95:PRO:HB2	1:A:229:TRP:CH2	2.51	0.46
2:B:244:ILE:HD12	2:B:244:ILE:H	1.80	0.46
2:B:297:GLU:HG3	2:B:298:GLU:OE2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:ARG:O	1:A:203:GLU:HG2	2.15	0.46
1:A:221:HIS:CD2	1:A:228:LEU:H	2.34	0.46
2:B:87:PHE:CE2	2:B:158:ALA:HB3	2.51	0.46
3:P:806:DT:H2'	3:P:807:DC:C6	2.51	0.46
2:B:66:LYS:HA	2:B:407:GLN:HE22	1.81	0.46
3:P:815:DG:H2'	3:P:816:DG:C8	2.51	0.46
1:A:94:ILE:HG23	1:A:230:MET:CE	2.46	0.45
1:A:116:PHE:HA	1:A:148:VAL:HG21	1.98	0.45
4:T:719:DG:H2'	4:T:720:DG:C8	2.51	0.45
1:A:96:HIS:NE2	1:A:269:GLN:NE2	2.64	0.45
1:A:441:TYR:HB2	1:A:458:VAL:HG12	1.97	0.45
2:B:61:PHE:N	2:B:61:PHE:CD1	2.84	0.45
2:B:289:LEU:HD12	2:B:289:LEU:HA	1.60	0.45
1:A:41:MET:CE	1:A:73:LYS:HD3	2.46	0.45
1:A:76:ASP:OD2	1:A:78:ARG:NH2	2.49	0.45
1:A:177:ASP:HB2	7:A:564:HOH:O	2.15	0.45
1:A:394:GLN:O	1:A:395:LYS:C	2.60	0.45
1:A:403:THR:CG2	1:A:404:GLU:N	2.74	0.45
1:A:281:LYS:C	1:A:283:LEU:H	2.24	0.45
1:A:339:TYR:CD1	1:A:375:ILE:HD11	2.52	0.45
1:A:493:VAL:HG22	1:A:494:ASN:N	2.32	0.45
1:A:84:THR:HG22	1:A:85:GLN:O	2.17	0.45
2:B:93:GLY:O	2:B:95:PRO:HD3	2.16	0.45
2:B:172:ARG:CZ	2:B:180:ILE:HB	2.46	0.45
2:B:173:LYS:O	2:B:176:PRO:HD3	2.17	0.45
2:B:13:LYS:NZ	2:B:88:TRP:HH2	2.11	0.45
3:P:810:DT:H2'	3:P:811:DG:H8	1.82	0.45
4:T:705:DT:H2''	4:T:706:DT:H5'	1.98	0.45
1:A:13:LYS:HG3	1:A:84:THR:O	2.17	0.45
1:A:356:ARG:CD	1:A:358:ARG:HG3	2.43	0.45
1:A:483:TYR:CE1	1:A:520:GLN:HB3	2.51	0.45
2:B:88:TRP:CE3	2:B:154:LYS:HE2	2.52	0.45
1:A:17:ASP:O	1:A:83:ARG:NH1	2.50	0.45
1:A:195:ILE:O	1:A:196:GLY:C	2.58	0.45
2:B:94:ILE:HG21	2:B:181:TYR:HE1	1.82	0.45
1:A:389:PHE:HB3	1:A:391:LEU:HD11	1.98	0.44
1:A:408:ALA:HB2	2:B:337:TRP:HH2	1.82	0.44
2:B:8:VAL:HA	2:B:9:PRO:HD3	1.73	0.44
2:B:157:PRO:HG3	2:B:184:MET:HA	1.99	0.44
2:B:246:LEU:HD11	2:B:264:LEU:HD21	1.99	0.44
4:T:706:DT:H2'	4:T:707:DG:H8	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:710:DG:H2'	4:T:711:DC:C6	2.52	0.44
1:A:358:ARG:HH11	1:A:358:ARG:CB	2.28	0.44
1:A:542:ILE:HG23	2:B:283:LEU:HD13	1.99	0.44
4:T:713:DC:C2'	4:T:714:DG:O4'	2.64	0.44
1:A:181:TYR:CD2	2:B:138:GLU:HA	2.52	0.44
1:A:408:ALA:HB2	2:B:337:TRP:CH2	2.52	0.44
1:A:410:TRP:CE2	2:B:363:ASN:ND2	2.86	0.44
1:A:458:VAL:CG1	1:A:548:VAL:HG22	2.46	0.44
2:B:30:LYS:O	2:B:34:LEU:HG	2.18	0.44
2:B:281:LYS:CB	2:B:281:LYS:NZ	2.80	0.44
1:A:381:VAL:HG22	2:B:25:PRO:HB3	2.00	0.44
2:B:63:ILE:HD11	2:B:408:ALA:O	2.17	0.44
2:B:115:TYR:O	2:B:149:LEU:HB2	2.17	0.44
1:A:404:GLU:OE2	1:A:509:GLN:NE2	2.51	0.44
2:B:150:PRO:HG2	2:B:153:TRP:CB	2.48	0.44
2:B:341:ILE:CG2	2:B:342:TYR:N	2.80	0.44
1:A:285:GLY:O	1:A:286:THR:C	2.60	0.44
1:A:450:THR:O	1:A:452:LEU:HD13	2.17	0.44
2:B:168:LEU:HB3	2:B:172:ARG:HG3	2.00	0.44
2:B:337:TRP:NE1	2:B:367:GLN:HG2	2.30	0.44
3:P:807:DC:H2'	3:P:808:DC:O4'	2.17	0.44
1:A:3:SER:HA	1:A:212:TRP:O	2.17	0.44
1:A:122:GLU:O	1:A:125:ARG:HB2	2.18	0.44
1:A:232:TYR:HB3	1:A:240:THR:O	2.18	0.44
1:A:494:ASN:HB3	2:B:289:LEU:HD22	2.00	0.44
1:A:500:GLN:HE21	2:B:420:PRO:HB2	1.81	0.44
1:A:302:GLU:O	1:A:305:GLU:HB2	2.18	0.44
1:A:312:GLU:HA	1:A:312:GLU:OE1	2.16	0.44
1:A:454:LYS:HB2	1:A:552:VAL:O	2.17	0.44
1:A:454:LYS:HE2	1:A:468:THR:HG22	1.99	0.44
1:A:525:LEU:HA	1:A:525:LEU:HD23	1.77	0.44
2:B:248:GLU:HB3	2:B:307:ARG:HH12	1.83	0.44
1:A:3:SER:HA	1:A:4:PRO:HD3	1.82	0.44
1:A:362:THR:HA	1:A:366:LYS:HE3	1.99	0.43
3:P:811:DG:H2''	3:P:812:DT:O5'	2.18	0.43
1:A:49:LYS:HB2	1:A:49:LYS:NZ	2.33	0.43
1:A:134:SER:C	1:A:136:ASN:H	2.26	0.43
1:A:337:TRP:CH2	1:A:368:LEU:HB2	2.53	0.43
2:B:100:LEU:HD13	2:B:179:VAL:HG13	1.99	0.43
2:B:163:SER:O	2:B:166:LYS:HB2	2.18	0.43
2:B:194:GLU:HB3	2:B:197:GLN:HG2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:ILE:HG22	1:A:119:PRO:HG3	1.97	0.43
2:B:175:ASN:N	2:B:176:PRO:HD3	2.33	0.43
2:B:260:LEU:HA	2:B:260:LEU:HD12	1.78	0.43
1:A:53:GLU:CD	1:A:53:GLU:N	2.65	0.43
1:A:244:ILE:HD13	1:A:267:ALA:HB2	2.01	0.43
1:A:399:GLU:O	1:A:403:THR:HB	2.19	0.43
1:A:56:TYR:O	1:A:143:ARG:NH2	2.51	0.43
1:A:254:VAL:HG13	1:A:289:LEU:HA	2.01	0.43
1:A:450:THR:O	1:A:451:LYS:HB2	2.19	0.43
4:T:722:DA:H2'	4:T:723:DC:H6	1.83	0.43
1:A:5:ILE:HG12	1:A:6:GLU:H	1.83	0.43
1:A:361:HIS:ND1	1:A:513:SER:CB	2.82	0.43
2:B:170:PRO:O	2:B:174:GLN:HG2	2.18	0.43
2:B:395:LYS:HG3	2:B:416:PHE:CE2	2.54	0.43
2:B:329:ILE:HG22	2:B:330:GLN:N	2.33	0.42
1:A:195:ILE:HG12	1:A:196:GLY:N	2.34	0.42
1:A:136:ASN:ND2	1:A:136:ASN:N	2.67	0.42
1:A:194:GLU:O	1:A:195:ILE:C	2.63	0.42
1:A:281:LYS:O	1:A:284:ARG:HG2	2.19	0.42
1:A:338:THR:HG22	1:A:339:TYR:N	2.34	0.42
1:A:443:ASP:OD2	1:A:498:ASP:OD2	2.36	0.42
2:B:78:ARG:HB2	2:B:78:ARG:HE	1.70	0.42
2:B:112:GLY:HA2	2:B:115:TYR:HD1	1.84	0.42
2:B:126:LYS:HE2	2:B:127:TYR:CZ	2.54	0.42
2:B:214:LEU:C	2:B:214:LEU:HD23	2.44	0.42
2:B:180:ILE:HG23	2:B:189:VAL:CG2	2.40	0.42
2:B:77:PHE:O	2:B:78:ARG:C	2.62	0.42
1:A:319:TYR:OH	1:A:385:LYS:HD3	2.19	0.42
1:A:541:GLY:HA2	1:A:546:GLU:HB2	2.01	0.42
2:B:16:MET:HE2	2:B:83:ARG:HA	2.02	0.42
2:B:26:LEU:HD12	2:B:133:PRO:HG2	1.98	0.42
2:B:54:ASN:HB3	2:B:143:ARG:HH12	1.84	0.42
2:B:81:ASN:OD1	2:B:153:TRP:HD1	2.03	0.42
2:B:377:THR:O	2:B:378:GLU:C	2.63	0.42
1:A:111:VAL:HB	1:A:185:ASP:HB2	2.02	0.42
1:A:169:GLU:O	1:A:170:PRO:C	2.62	0.42
2:B:271:TYR:HA	2:B:272:PRO:HD3	1.81	0.42
2:B:303:LEU:HA	2:B:306:ASN:HB2	2.01	0.42
2:B:422:LEU:H	2:B:422:LEU:CD2	2.31	0.42
1:A:524:GLN:HA	1:A:524:GLN:OE1	2.19	0.42
2:B:120:LEU:O	2:B:121:ASP:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:170:PRO:O	2:B:171:PHE:C	2.63	0.42
2:B:237:ASP:C	2:B:238:LYS:HG3	2.45	0.42
4:T:704:DG:N2	4:T:704:DG:OP2	2.51	0.42
2:B:61:PHE:CD2	2:B:403:THR:HB	2.55	0.42
2:B:112:GLY:HA2	2:B:115:TYR:CD1	2.54	0.42
2:B:341:ILE:HG22	2:B:342:TYR:N	2.35	0.42
1:A:88:TRP:CE2	2:B:57:ASN:HB2	2.55	0.41
1:A:91:GLN:O	4:T:708:DG:H4'	2.20	0.41
1:A:216:THR:C	1:A:217:PRO:O	2.62	0.41
1:A:222:GLN:O	1:A:224:GLU:N	2.53	0.41
1:A:277:ARG:HH12	1:A:357:MET:HB2	1.84	0.41
1:A:288:ALA:O	1:A:291:GLU:HB3	2.20	0.41
1:A:30:LYS:HG2	1:A:71:TRP:HZ3	1.80	0.41
1:A:270:ILE:HD12	1:A:270:ILE:HA	1.82	0.41
1:A:290:THR:O	1:A:291:GLU:C	2.63	0.41
1:A:305:GLU:O	1:A:309:ILE:HG13	2.20	0.41
1:A:457:TYR:CD1	1:A:457:TYR:C	2.97	0.41
2:B:236:PRO:O	2:B:238:LYS:N	2.51	0.41
2:B:320:ASP:HB3	2:B:323:LYS:HB2	2.02	0.41
2:B:391:LEU:HA	2:B:392:PRO:HD3	1.86	0.41
1:A:266:TRP:CE2	3:P:820:DC:H4'	2.55	0.41
2:B:80:LEU:HD12	2:B:80:LEU:O	2.19	0.41
2:B:164:MET:SD	2:B:167:ILE:HD12	2.60	0.41
1:A:34:LEU:CD2	1:A:62:ALA:HB2	2.43	0.41
1:A:513:SER:O	1:A:519:ASN:ND2	2.53	0.41
2:B:10:VAL:HG13	2:B:85:GLN:NE2	2.35	0.41
2:B:266:TRP:C	2:B:268:SER:N	2.76	0.41
2:B:332:GLN:CG	2:B:338:THR:HG23	2.50	0.41
1:A:358:ARG:HB3	1:A:358:ARG:NH1	2.27	0.41
1:A:412:PRO:O	1:A:414:TRP:HD1	2.03	0.41
1:A:540:LYS:HA	1:A:540:LYS:HD3	1.70	0.41
2:B:422:LEU:N	2:B:422:LEU:CD2	2.84	0.41
4:T:722:DA:H2''	4:T:723:DC:O5'	2.20	0.41
1:A:24:TRP:CH2	4:T:704:DG:C8	3.08	0.41
1:A:219:LYS:H	1:A:219:LYS:HG2	1.73	0.41
2:B:209:LEU:HG	2:B:210:LEU:N	2.36	0.41
1:A:31:ILE:CD1	1:A:135:ILE:HA	2.50	0.41
1:A:282:LEU:HD21	1:A:296:THR:HB	2.02	0.41
1:A:293:ILE:O	1:A:293:ILE:HG23	2.19	0.41
1:A:472:THR:OG1	1:A:473:THR:N	2.53	0.41
1:A:489:SER:HB2	1:A:493:VAL:HB	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:92:LEU:CD2	2:B:94:ILE:HG13	2.49	0.41
2:B:331:LYS:HA	2:B:337:TRP:CE3	2.55	0.41
1:A:100:LEU:O	1:A:318:TYR:HB3	2.20	0.41
1:A:105:SER:HB3	1:A:198:HIS:CG	2.56	0.41
1:A:543:GLY:HA2	1:A:544:GLY:HA2	1.41	0.41
2:B:180:ILE:HG12	2:B:189:VAL:CG1	2.50	0.41
1:A:41:MET:HE1	1:A:73:LYS:HD3	2.02	0.41
1:A:134:SER:C	1:A:136:ASN:N	2.79	0.41
1:A:419:THR:HG23	1:A:419:THR:O	2.21	0.41
2:B:8:VAL:HG13	2:B:8:VAL:O	2.21	0.41
1:A:221:HIS:ND1	1:A:221:HIS:N	2.69	0.41
1:A:416:PHE:HZ	1:A:422:LEU:HD11	1.85	0.41
2:B:5:ILE:HG12	2:B:6:GLU:N	2.36	0.41
2:B:300:GLU:O	2:B:304:ALA:HB2	2.21	0.41
1:A:239:TRP:HE3	1:A:240:THR:N	2.18	0.40
1:A:345:PRO:HB2	1:A:346:PHE:HD1	1.86	0.40
1:A:424:LYS:HB3	1:A:424:LYS:HE3	1.78	0.40
2:B:74:LEU:HD12	2:B:74:LEU:HA	1.96	0.40
2:B:246:LEU:HD22	2:B:260:LEU:HD11	2.03	0.40
2:B:337:TRP:HE1	2:B:367:GLN:CG	2.33	0.40
2:B:162:SER:OG	2:B:163:SER:N	2.54	0.40
2:B:207:GLN:C	2:B:209:LEU:H	2.29	0.40
1:A:149:LEU:HA	1:A:150:PRO:HD3	1.87	0.40
1:A:205:LEU:HD22	1:A:209:LEU:CD2	2.51	0.40
1:A:287:LYS:HE3	1:A:289:LEU:HG	2.01	0.40
2:B:148:VAL:O	2:B:149:LEU:C	2.64	0.40
2:B:242:GLN:HA	2:B:243:PRO:HD3	1.79	0.40
2:B:263:LYS:HE3	2:B:263:LYS:HB2	1.79	0.40
1:A:337:TRP:CZ3	1:A:368:LEU:HB2	2.57	0.40
1:A:491:LEU:HG	1:A:529:GLU:CG	2.51	0.40
1:A:540:LYS:HB3	1:A:542:ILE:CD1	2.52	0.40
2:B:341:ILE:N	2:B:341:ILE:CD1	2.83	0.40
1:A:134:SER:HB3	1:A:139:THR:HB	2.02	0.40
1:A:407:GLN:O	1:A:408:ALA:C	2.63	0.40
2:B:302:GLU:HG2	2:B:306:ASN:ND2	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:173:LYS:NZ	2:B:173:LYS:NZ[3_656]	2.01	0.19

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	553/560 (99%)	501 (91%)	44 (8%)	8 (1%)	9	30
2	B	401/452 (89%)	355 (88%)	37 (9%)	9 (2%)	5	20
All	All	954/1012 (94%)	856 (90%)	81 (8%)	17 (2%)	6	25

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	LYS
2	B	5	ILE
2	B	193	LEU
1	A	217	PRO
1	A	291	GLU
2	B	240	THR
2	B	357	MET
1	A	223	LYS
1	A	284	ARG
2	B	170	PRO
2	B	14	PRO
2	B	87	PHE
1	A	419	THR
2	B	234	LEU
2	B	237	ASP
1	A	157	PRO
1	A	490	GLY

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	495/500 (99%)	446 (90%)	49 (10%)	7	24
2	B	369/411 (90%)	331 (90%)	38 (10%)	7	23
All	All	864/911 (95%)	777 (90%)	87 (10%)	7	24

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	TRP
1	A	42	GLU
1	A	49	LYS
1	A	92	LEU
1	A	100	LEU
1	A	109	LEU
1	A	135	ILE
1	A	136	ASN
1	A	137	ASN
1	A	168	LEU
1	A	195	ILE
1	A	201	LYS
1	A	205	LEU
1	A	209	LEU
1	A	216	THR
1	A	218	ASP
1	A	221	HIS
1	A	228	LEU
1	A	254	VAL
1	A	264	LEU
1	A	268	SER
1	A	270	ILE
1	A	284	ARG
1	A	290	THR
1	A	295	LEU
1	A	300	GLU
1	A	302	GLU
1	A	312	GLU
1	A	328	GLU
1	A	330	GLN
1	A	331	LYS
1	A	340	GLN
1	A	356	ARG
1	A	358	ARG

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Mol	Chain	Res	Type
1	A	373	GLN
1	A	391	LEU
1	A	393	ILE
1	A	394	GLN
1	A	399	GLU
1	A	402	TRP
1	A	403	THR
1	A	415	GLU
1	A	458	VAL
1	A	469	LEU
1	A	480	GLN
1	A	491	LEU
1	A	507	GLN
1	A	523	GLU
1	A	548	VAL
2	B	6	GLU
2	B	16	MET
2	B	35	VAL
2	B	44	GLU
2	B	53	GLU
2	B	61	PHE
2	B	66	LYS
2	B	70	LYS
2	B	72	ARG
2	B	78	ARG
2	B	91	GLN
2	B	92	LEU
2	B	122	GLU
2	B	135	ILE
2	B	148	VAL
2	B	194	GLU
2	B	232	TYR
2	B	242	GLN
2	B	268	SER
2	B	277	ARG
2	B	281	LYS
2	B	284	ARG
2	B	286	THR
2	B	289	LEU
2	B	295	LEU
2	B	297	GLU
2	B	303	LEU

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Mol	Chain	Res	Type
2	B	314	VAL
2	B	324	ASP
2	B	325	LEU
2	B	336	GLN
2	B	364	ASP
2	B	368	LEU
2	B	372	VAL
2	B	374	LYS
2	B	396	GLU
2	B	418	ASN
2	B	428	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	ASN
1	A	137	ASN
1	A	161	GLN
1	A	197	GLN
1	A	242	GLN
1	A	255	ASN
1	A	269	GLN
1	A	306	ASN
1	A	336	GLN
1	A	373	GLN
1	A	464	GLN
1	A	480	GLN
1	A	487	GLN
1	A	500	GLN
1	A	509	GLN
1	A	545	ASN
1	A	547	GLN
2	B	91	GLN
2	B	182	GLN
2	B	235	HIS
2	B	242	GLN
2	B	265	ASN
2	B	334	GLN
2	B	336	GLN
2	B	367	GLN
2	B	394	GLN
2	B	428	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	URT	P	822	3	16,23,24	1.64	4 (25%)	19,33,36	0.98	0

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	URT	P	822	3	-	3/5/21/22	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	822	URT	C6-C5	3.27	1.44	1.39
3	P	822	URT	O2-C1	-2.88	1.36	1.42
3	P	822	URT	C5-N3	2.57	1.39	1.34
3	P	822	URT	C8-N4	2.27	1.38	1.33

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	P	822	URT	O1-C1-O2-C10
3	P	822	URT	C3-C4-N1-C7

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Mol	Chain	Res	Type	Atoms
3	P	822	URT	C3-C4-N1-C5

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
6	SO4	T	2	-	4,4,4	0.37	0	6,6,6	0.13	0
6	SO4	P	3	-	4,4,4	0.39	0	6,6,6	0.17	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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 [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	A	555/560 (99%)	-0.38	3 (0%) 87 83	30, 59, 110, 135	0
2	B	407/452 (90%)	-0.19	2 (0%) 87 83	40, 78, 133, 145	0
3	P	17/21 (80%)	-0.62	0 100 100	37, 65, 84, 95	0
4	T	23/27 (85%)	-0.27	0 100 100	36, 76, 125, 160	0
All	All	1002/1060 (94%)	-0.31	5 (0%) 87 83	30, 67, 124, 160	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	555	GLY	3.5
2	B	214	LEU	2.7
2	B	430	GLU	2.2
1	A	1	PRO	2.2
1	A	419	THR	2.1

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(Å ²)	Q<0.9
3	URT	P	822	21/22	0.95	0.06	45,49,51,52	0

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
6	SO4	P	3	5/5	0.78	0.10	110,111,111,112	0
6	SO4	T	2	5/5	0.94	0.12	102,102,103,103	0
5	MG	A	601	1/1	0.99	0.12	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.