



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:52 PM UTC

PDB ID : 6P71 / pdb_00006p71
Title : X-ray crystal structure of a bacterial reiterative transcription complex of pyrBI promoter
Authors : Shin, Y.; Murakami, K.S.
Deposited on : 2019-06-04
Resolution : 2.92 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

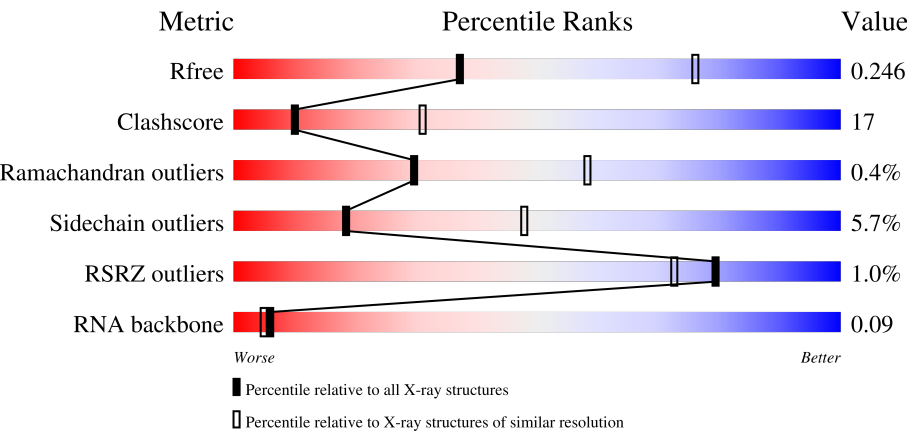
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.92 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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)
RNA backbone	3983	1016 (3.12-2.72)

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	315	45% 24% • 28%
1	B	315	43% 26% • 30%
2	C	1119	% 64% 32% ••
3	D	1524	% 67% 28% ••

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Mol	Chain	Length	Quality of chain
4	E	99	3% 76% 18% • 5%
5	F	423	% 55% 25% • 18%
6	G	21	5% 29% 38% 5% 29%
7	H	27	26% 33% 41%
8	I	4	25% 75%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 28417 atoms, of which 11 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 DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	0	0
			1782	1138	310	332	2			
1	B	222	Total	C	N	O	S	0	0	0
			1750	1118	304	326	2			

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	1111	Total	C	N	O	S	0	0	0
			8747	5536	1560	1627	24			

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	1500	Total	C	N	O	S	0	0	0
			11820	7489	2083	2212	36			

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	94	Total	C	N	O	S	0	0	0
			761	486	132	139	4			

- Molecule 5 is a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	346	Total	C	N	O	S	0	0	0
			2796	1764	508	520	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	46	THR	ALA	conflict	UNP Q72L95

- Molecule 6 is a DNA chain called DNA (5'-D(P*CP*TP*CP*CP*CP*GP*GP*CP*AP*AP*AP*TP*TP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	G	15	Total	C	N	O	P	0	0	0
			304	145	53	91	15			

- Molecule 7 is a DNA chain called DNA (5'-D(*TP*AP*TP*AP*AP*TP*CP*GP*AP*GP*CP*CP*GP*GP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	H	16	Total	C	N	O	P	0	0	0
			329	157	65	92	15			

- Molecule 8 is a RNA chain called RNA (5'-R(P*AP*AP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	I	4	Total	C	N	O	P	0	0	0
			84	38	14	28	4			

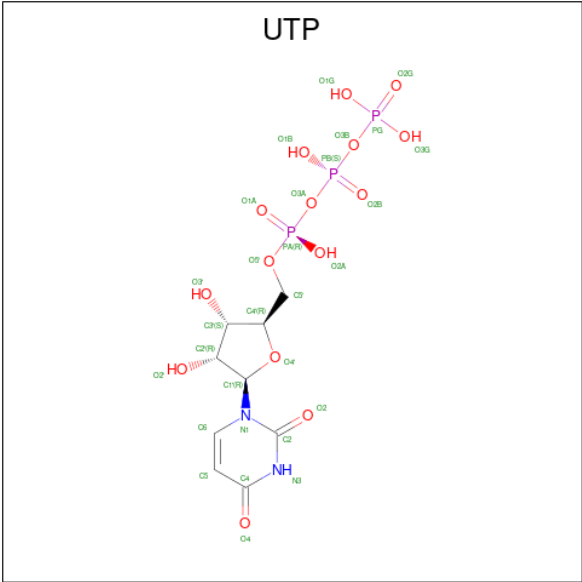
- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	D	2	Total	Mg	0	0
			2	2		

- Molecule 10 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	2	Total	Zn	0	0
			2	2		

- Molecule 11 is URIDINE 5'-TRIPHOSPHATE (CCD ID: UTP) (formula: C₉H₁₅N₂O₁₅P₃).

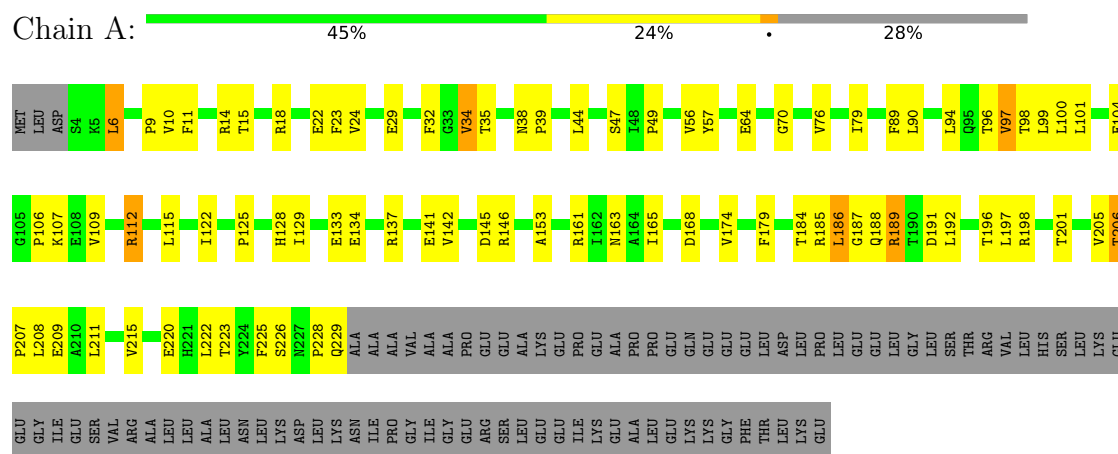


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
11	D	1	Total 40	C 9	H 11	N 2	O 15	P 3	0	0

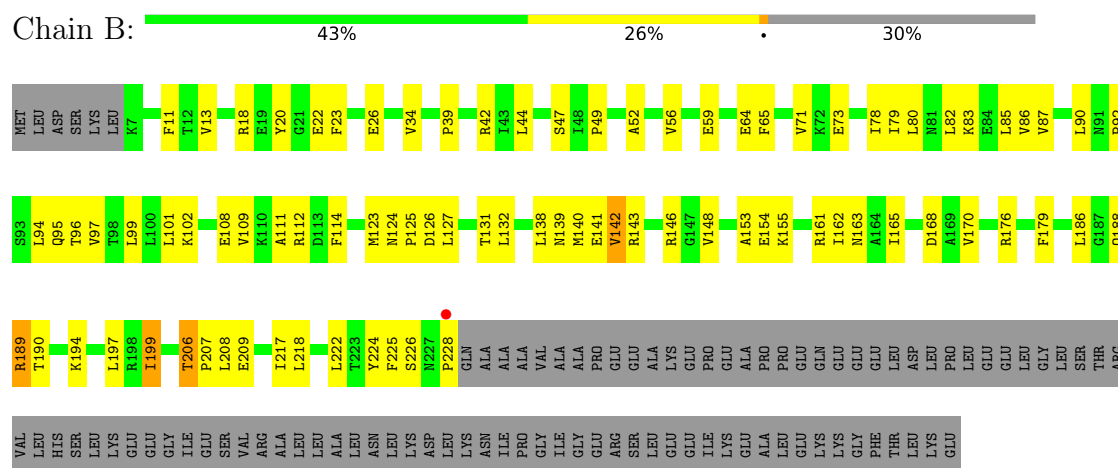
3 Residue-property plots

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.

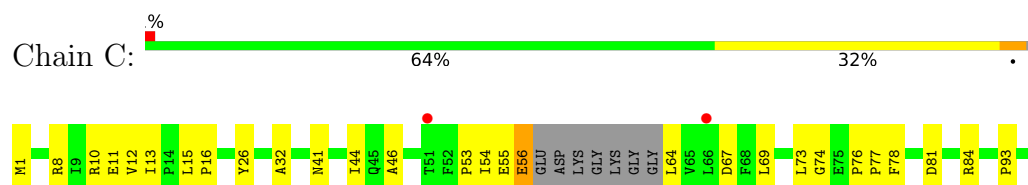
• Molecule 1: DNA-directed RNA polymerase subunit alpha

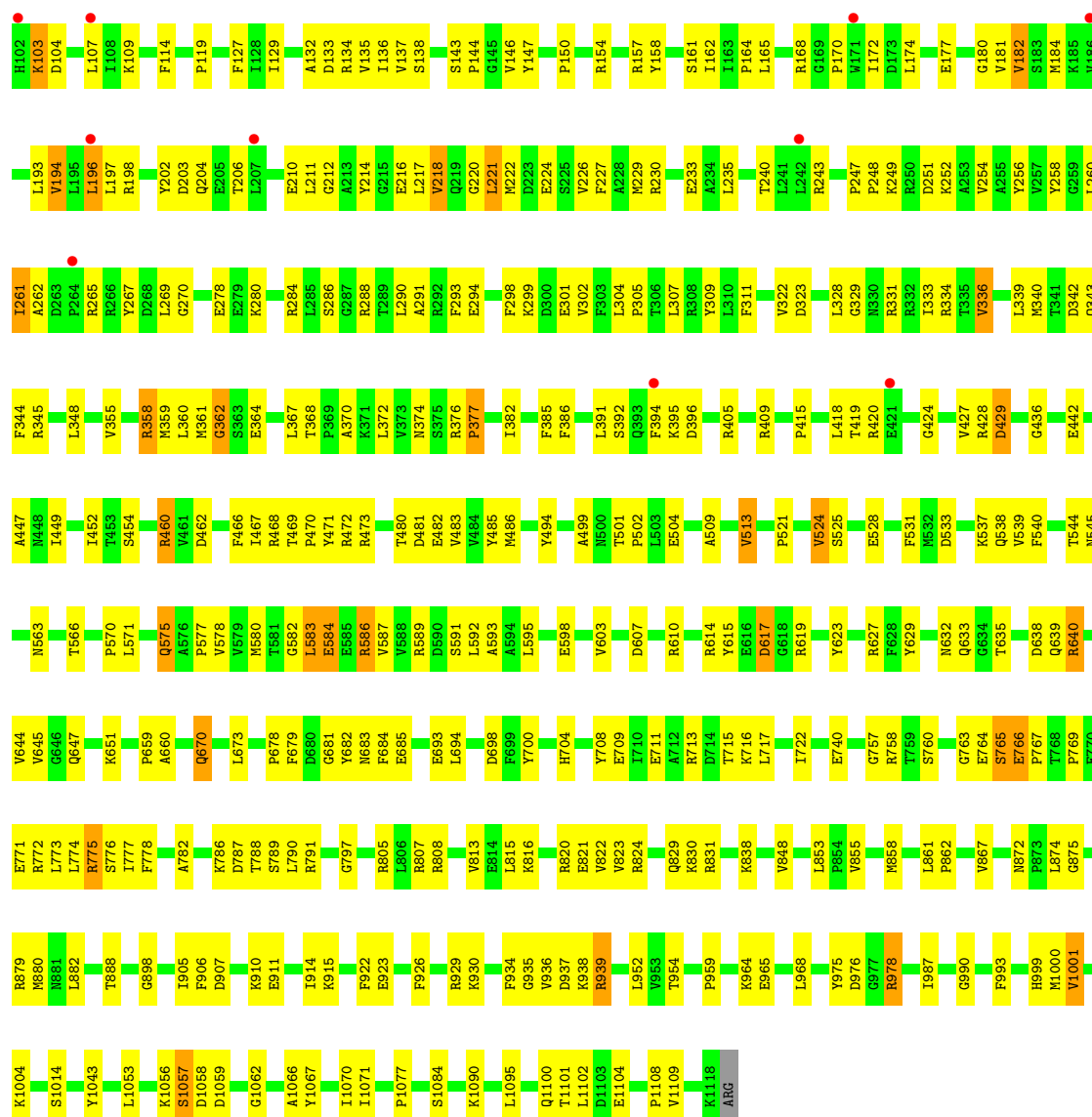


• Molecule 1: DNA-directed RNA polymerase subunit alpha

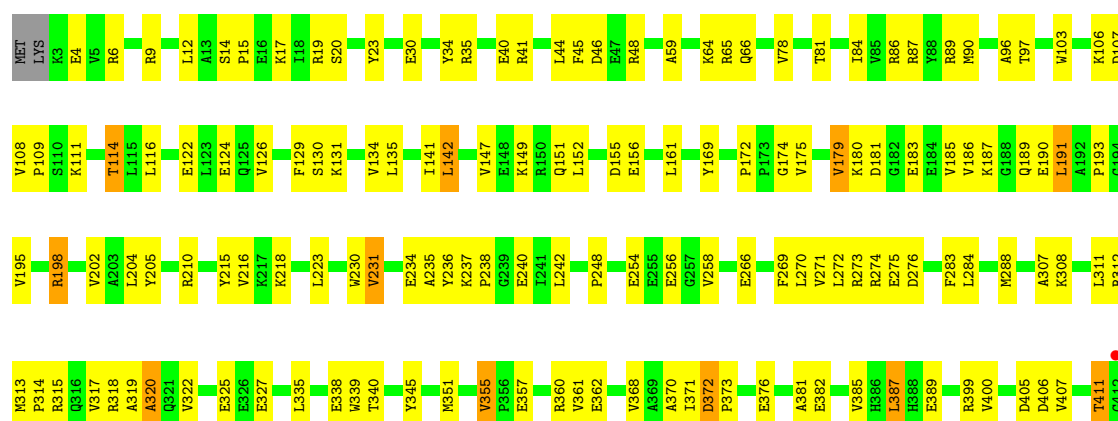


• Molecule 2: DNA-directed RNA polymerase subunit beta



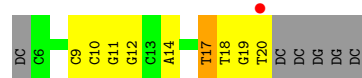


• Molecule 3: DNA-directed RNA polymerase subunit beta'

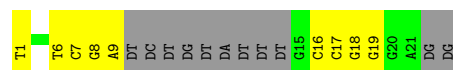
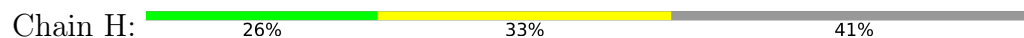




- Molecule 6: DNA (5'-D(P*CP*TP*CP*CP*CP*GP*GP*CP*AP*AP*AP*TP*TP*GP*T)-3')



- Molecule 7: DNA (5'-D(*TP*AP*TP*AP*AP*TP*CP*GP*AP*GP*CP*CP*GP*GP*GP*A)-3')



- Molecule 8: RNA (5'-R(P*AP*AP*UP*U)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	187.51Å 100.66Å 296.30Å 90.00° 98.32° 90.00°	Depositor
Resolution (Å)	48.97 – 2.92 48.97 – 2.92	Depositor EDS
% Data completeness (in resolution range)	98.1 (48.97-2.92) 98.0 (48.97-2.92)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.91Å)	Xtriage
Refinement program	PHENIX (1.14_3260)	Depositor
R, R_{free}	0.211 , 0.248 0.211 , 0.246	Depositor DCC
R_{free} test set	1970 reflections (1.58%)	wwPDB-VP
Wilson B-factor (Å ²)	75.8	Xtriage
Anisotropy	0.540	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	28417	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ZN, UTP

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	0.54	0/1814	0.89	0/2466
1	B	0.56	0/1782	0.90	0/2424
2	C	0.55	0/8914	0.88	5/12060 (0.0%)
3	D	0.59	0/12028	0.91	6/16265 (0.0%)
4	E	0.52	0/775	0.84	0/1045
5	F	0.50	0/2841	0.85	0/3824
6	G	0.61	1/339 (0.3%)	0.62	0/520
7	H	0.43	0/369	0.55	0/566
8	I	1.16	2/93 (2.2%)	1.03	1/142 (0.7%)
All	All	0.56	3/28955 (0.0%)	0.89	12/39312 (0.0%)

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
3	D	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	I	1	A	O3'-P	-6.09	1.52	1.61
8	I	2	U	O3'-P	-5.25	1.53	1.61
6	G	17	DT	O3'-P	-5.13	1.53	1.61

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1235	GLN	CA-CB-CG	-7.59	98.92	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	782	ALA	N-CA-C	6.72	120.12	110.24
8	I	0	A	C2'-C3'-O3'	-6.65	103.72	113.70
2	C	682	TYR	CA-C-N	5.79	132.87	122.38
2	C	682	TYR	C-N-CA	5.79	132.87	122.38
3	D	1455	LYS	CD-CE-NZ	5.78	130.39	111.90
3	D	254	GLU	CA-CB-CG	-5.77	102.57	114.10
3	D	740	PHE	N-CA-C	5.20	118.44	112.57
3	D	1344	VAL	CG1-CB-CG2	-5.15	99.47	110.80
2	C	765	SER	CA-C-N	5.04	134.11	121.80
2	C	765	SER	C-N-CA	5.04	134.11	121.80
3	D	1208	ASP	N-CA-C	-5.02	100.32	108.41

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	D	583	ASP	Mainchain

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	1782	0	1834	73	0
1	B	1750	0	1797	83	0
2	C	8747	0	8837	346	0
3	D	11820	0	12038	424	2
4	E	761	0	778	23	0
5	F	2796	0	2865	101	0
6	G	304	0	170	29	0
7	H	329	0	182	13	0
8	I	84	0	42	16	0
9	D	2	0	0	0	0
10	D	2	0	0	0	0
11	D	29	11	11	1	0
All	All	28406	11	28554	986	2

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 17.

All (986) 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)
6:G:17:DT:H3	8:I:1:A:N6	1.07	1.49
6:G:18:DT:H3	8:I:0:A:N6	1.25	1.32
6:G:17:DT:O2	8:I:1:A:N1	1.78	1.15
2:C:853:LEU:HB2	2:C:858:MET:HE1	1.31	1.10
3:D:218:LYS:HG2	3:D:338:GLU:HG2	1.34	1.09
3:D:949:ILE:HD11	3:D:1023:MET:HE1	1.32	1.09
3:D:1493:LYS:HA	3:D:1493:LYS:HE3	1.46	0.95
2:C:221:LEU:HD11	2:C:307:LEU:HD21	1.47	0.95
6:G:18:DT:O2	8:I:0:A:N1	2.00	0.94
5:F:168:LYS:HG3	5:F:172:ARG:HH21	1.33	0.93
2:C:617:ASP:OD2	2:C:619:ARG:NE	2.01	0.93
3:D:284:LEU:HD22	3:D:288:MET:HE2	1.51	0.93
3:D:41:ARG:HA	3:D:46:ASP:OD2	1.69	0.92
1:B:138:LEU:HD21	1:B:140:MET:HE3	1.49	0.92
3:D:566:ILE:HD11	5:F:192:LEU:HD21	1.48	0.92
3:D:664:LYS:NZ	3:D:693:GLU:OE1	2.03	0.92
2:C:109:LYS:HG2	2:C:368:THR:HG22	1.48	0.92
5:F:193:ARG:HB3	7:H:7:DC:H5"	1.51	0.92
1:B:108:GLU:HG2	1:B:131:THR:HG22	1.52	0.91
2:C:55:GLU:O	2:C:56:GLU:HB3	1.71	0.91
2:C:615:TYR:HH	2:C:623:TYR:HH	1.13	0.90
3:D:1154:GLU:HG2	3:D:1159:ARG:HA	1.54	0.90
5:F:338:LEU:HD23	5:F:339:PRO:HD2	1.54	0.89
2:C:212:GLY:HA2	2:C:218:VAL:HG21	1.54	0.89
3:D:96:ALA:HB3	3:D:554:LEU:HD23	1.54	0.88
6:G:17:DT:C2	8:I:1:A:N1	2.40	0.88
2:C:853:LEU:HB2	2:C:858:MET:CE	2.03	0.88
1:B:206:THR:HG22	1:B:209:GLU:H	1.37	0.88
2:C:32:ALA:HB2	2:C:73:LEU:HD12	1.56	0.88
3:D:1143:GLY:O	3:D:1147:ARG:HD2	1.74	0.88
1:B:176:ARG:NH2	3:D:888:GLU:OE1	2.06	0.87
3:D:238:PRO:HD3	3:D:318:ARG:HG3	1.56	0.87
2:C:229:MET:HB2	2:C:233:GLU:HB2	1.55	0.86
3:D:65:ARG:HB2	3:D:65:ARG:HH11	1.39	0.85
2:C:394:PHE:HE2	2:C:632:ASN:HB3	1.40	0.85
2:C:168:ARG:NH1	2:C:345:ARG:HD3	1.92	0.85
3:D:322:VAL:HG22	3:D:335:LEU:CD2	2.08	0.84
2:C:56:GLU:HB2	2:C:359:MET:HE3	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:I:0:A:O2'	8:I:1:A:O5'	1.96	0.84
2:C:1070:ILE:HG21	3:D:655:PRO:HB2	1.60	0.83
2:C:358:ARG:HB3	2:C:372:LEU:HD12	1.60	0.83
3:D:1486:VAL:HG21	4:E:22:VAL:HG13	1.60	0.83
3:D:351:MET:HE2	3:D:368:VAL:HG11	1.61	0.83
2:C:146:VAL:HG22	2:C:162:ILE:HG12	1.60	0.83
2:C:333:ILE:HD11	2:C:467:ILE:HD11	1.61	0.83
1:A:188:GLN:HG2	1:A:189:ARG:HG2	1.59	0.82
3:D:1496:GLU:OE1	3:D:1500:LYS:NZ	2.13	0.82
2:C:521:PRO:HB3	3:D:1068:LEU:HD21	1.61	0.81
1:A:222:LEU:HD21	1:B:218:LEU:HD23	1.62	0.81
6:G:18:DT:H2'	6:G:19:DG:C8	2.16	0.81
2:C:294:GLU:HB3	2:C:299:LYS:HD2	1.63	0.80
3:D:566:ILE:HD11	5:F:192:LEU:CD2	2.11	0.80
3:D:963:TYR:CE2	3:D:1002:LYS:HD3	2.16	0.80
3:D:1263:PHE:HA	3:D:1375:MET:HE1	1.64	0.80
2:C:16:PRO:HB3	2:C:460:ARG:NH1	1.97	0.79
2:C:286:SER:OG	2:C:301:GLU:OE2	1.99	0.79
2:C:374:ASN:OD1	5:F:276:ARG:HD2	1.81	0.79
3:D:1486:VAL:CG2	4:E:22:VAL:HG13	2.12	0.79
1:B:49:PRO:HA	1:B:148:VAL:HG12	1.66	0.78
2:C:343:GLN:HG3	2:C:385:PHE:HB2	1.65	0.78
2:C:615:TYR:OH	2:C:623:TYR:OH	1.97	0.78
2:C:358:ARG:HB3	2:C:372:LEU:CD1	2.14	0.78
3:D:14:SER:HB3	3:D:511:TRP:CE2	2.19	0.78
3:D:1283:ILE:HG13	3:D:1315:ASP:CG	2.09	0.78
2:C:180:GLY:O	2:C:220:GLY:HA3	1.85	0.77
2:C:766:GLU:CB	2:C:767:PRO:HD2	2.15	0.77
3:D:65:ARG:HB2	3:D:65:ARG:NH1	1.99	0.77
3:D:114:THR:HG21	3:D:498:VAL:HG21	1.66	0.77
2:C:470:PRO:HD3	2:C:485:TYR:CE2	2.20	0.77
2:C:524:VAL:HG13	2:C:528:GLU:HB2	1.67	0.77
3:D:1283:ILE:HG13	3:D:1315:ASP:OD2	1.84	0.77
5:F:400:ILE:HG22	5:F:403:LYS:HE2	1.66	0.77
6:G:18:DT:H2''	6:G:19:DG:H5'	1.66	0.77
1:A:153:ALA:N	1:A:168:ASP:OD1	2.14	0.76
3:D:124:GLU:OE2	3:D:587:ARG:NH2	2.19	0.76
1:A:56:VAL:HG22	1:A:142:VAL:HG12	1.66	0.76
2:C:504:GLU:HG2	2:C:509:ALA:HB2	1.67	0.76
3:D:437:VAL:HG11	5:F:175:HIS:CD2	2.20	0.76
2:C:769:PRO:HG3	3:D:65:ARG:NH2	2.01	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1084:THR:HG23	3:D:1238:MET:HA	1.67	0.76
2:C:607:ASP:HB2	2:C:610:ARG:NH1	2.00	0.75
2:C:711:GLU:HG2	2:C:822:VAL:HG22	1.68	0.75
3:D:1084:THR:O	3:D:1088:THR:HG23	1.87	0.75
5:F:153:PRO:HA	5:F:156:VAL:HG22	1.69	0.74
3:D:314:PRO:HB2	3:D:317:VAL:HG12	1.69	0.74
1:B:80:LEU:HB3	3:D:867:ARG:NH2	2.03	0.74
5:F:168:LYS:HG3	5:F:172:ARG:NH2	2.02	0.74
2:C:470:PRO:HD3	2:C:485:TYR:HE2	1.53	0.74
3:D:904:VAL:HG22	3:D:905:PRO:HD2	1.70	0.73
3:D:1108:ARG:NH2	3:D:1198:TYR:O	2.21	0.73
2:C:164:PRO:HA	2:C:269:LEU:HD12	1.69	0.73
2:C:204:GLN:NE2	2:C:222:MET:SD	2.61	0.73
2:C:617:ASP:HB2	2:C:619:ARG:HG2	1.70	0.73
3:D:1386:ASP:CG	3:D:1413:THR:HG22	2.13	0.73
3:D:204:LEU:HD22	3:D:441:ARG:CZ	2.19	0.73
3:D:658:LEU:HD23	3:D:661:MET:CE	2.19	0.72
3:D:1047:LYS:HG2	3:D:1053:PHE:CE1	2.24	0.72
5:F:405:LEU:O	5:F:409:LYS:HG3	1.89	0.72
2:C:168:ARG:HH11	2:C:345:ARG:HD3	1.51	0.72
2:C:170:PRO:HD2	2:C:267:TYR:CE1	2.24	0.72
3:D:180:LYS:NZ	3:D:357:GLU:OE1	2.20	0.72
2:C:136:ILE:HB	2:C:336:VAL:HG13	1.72	0.72
2:C:764:GLU:C	2:C:766:GLU:H	1.98	0.72
2:C:936:VAL:HG11	2:C:959:PRO:HB2	1.71	0.72
6:G:11:DG:H2'	6:G:12:DG:H5''	1.72	0.72
2:C:704:HIS:CE1	2:C:1000:MET:HE1	2.24	0.71
3:D:500:ARG:NH1	3:D:1390:LEU:HD21	2.05	0.71
2:C:772:ARG:NH2	5:F:380:GLU:OE1	2.23	0.71
1:A:179:PHE:HB3	1:A:197:LEU:HD23	1.72	0.71
1:A:201:THR:HG21	1:A:205:VAL:O	1.90	0.71
3:D:959:GLU:HB3	3:D:963:TYR:CE1	2.26	0.71
2:C:394:PHE:CE2	2:C:632:ASN:HB3	2.24	0.71
3:D:1493:LYS:NZ	3:D:1496:GLU:HG3	2.06	0.71
5:F:130:VAL:HG21	5:F:159:ILE:HG21	1.73	0.70
3:D:1495:ILE:HG12	4:E:88:GLU:CG	2.21	0.70
3:D:1263:PHE:HD2	3:D:1375:MET:CE	2.05	0.70
2:C:181:VAL:HG23	2:C:220:GLY:O	1.91	0.70
5:F:289:GLU:O	5:F:293:GLU:HG3	1.92	0.70
3:D:704:ARG:HB2	3:D:745:MET:HG2	1.74	0.70
2:C:915:LYS:NZ	3:D:952:ASP:OD2	2.25	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:132:ALA:HB1	2:C:394:PHE:HE1	1.57	0.69
5:F:393:THR:HG22	5:F:395:GLU:H	1.56	0.69
2:C:937:ASP:OD1	2:C:939:ARG:HG2	1.92	0.69
3:D:1495:ILE:HG12	4:E:88:GLU:HG3	1.74	0.69
2:C:165:LEU:HD11	2:C:270:GLY:HA2	1.72	0.69
3:D:959:GLU:HB3	3:D:963:TYR:HE1	1.57	0.69
2:C:344:PHE:HD2	2:C:382:ILE:HD11	1.58	0.69
3:D:658:LEU:HD23	3:D:661:MET:HE1	1.74	0.69
6:G:12:DG:H5''	6:G:12:DG:H8	1.58	0.69
3:D:322:VAL:HG22	3:D:335:LEU:HD21	1.73	0.69
1:A:99:LEU:HB2	1:A:142:VAL:HG23	1.75	0.69
3:D:45:PHE:O	3:D:86:ARG:NH2	2.26	0.69
1:B:80:LEU:HD22	3:D:844:ALA:HA	1.75	0.68
3:D:657:LEU:HG	3:D:661:MET:HE2	1.74	0.68
5:F:115:LYS:HD3	5:F:173:TYR:CZ	2.28	0.68
2:C:41:ASN:OD1	2:C:46:ALA:HA	1.93	0.68
5:F:166:LEU:HD13	5:F:170:HIS:HB3	1.75	0.68
1:B:143:ARG:HG2	1:B:143:ARG:HH11	1.59	0.68
3:D:258:VAL:HG12	3:D:273:ARG:O	1.93	0.68
3:D:501:ALA:HB1	3:D:1452:ILE:HG22	1.75	0.68
2:C:74:GLY:HA3	2:C:93:PRO:HG2	1.75	0.68
2:C:340:MET:HE2	2:C:382:ILE:HD13	1.75	0.68
2:C:376:ARG:NH1	5:F:279:GLN:OE1	2.26	0.68
3:D:270:LEU:HD23	3:D:271:VAL:N	2.09	0.68
3:D:1364:HIS:ND1	3:D:1366:LYS:HB2	2.09	0.68
2:C:12:VAL:HG11	2:C:472:ARG:HD3	1.77	0.67
3:D:1087:ARG:CB	3:D:1237:THR:HG21	2.24	0.67
2:C:460:ARG:HD2	2:C:485:TYR:CZ	2.29	0.67
3:D:44:LEU:HB3	3:D:525:ARG:NH2	2.10	0.67
3:D:351:MET:HE2	3:D:368:VAL:CG1	2.25	0.67
3:D:1140:ILE:HG22	3:D:1144:LEU:HD12	1.76	0.67
1:A:97:VAL:HG13	1:A:98:THR:N	2.10	0.67
2:C:11:GLU:OE2	2:C:537:LYS:HE2	1.95	0.67
5:F:353:GLU:OE1	5:F:417:LYS:HD3	1.95	0.67
1:B:56:VAL:HG23	1:B:142:VAL:HG12	1.77	0.67
3:D:357:GLU:HB2	3:D:387:LEU:HD23	1.76	0.67
3:D:411:THR:HA	3:D:435:VAL:HG12	1.75	0.67
3:D:1234:THR:HG23	3:D:1238:MET:HE3	1.76	0.67
2:C:329:GLY:O	2:C:466:PHE:HE2	1.78	0.66
2:C:926:PHE:CZ	2:C:930:LYS:HD3	2.31	0.66
2:C:954:THR:HG23	2:C:965:GLU:OE2	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1102:LEU:HD23	2:C:1108:PRO:HA	1.77	0.66
3:D:116:LEU:HD21	3:D:465:LEU:HD23	1.77	0.66
6:G:17:DT:N3	8:I:1:A:N6	1.91	0.66
2:C:261:ILE:CG2	2:C:291:ALA:HB3	2.25	0.66
3:D:1147:ARG:HD3	3:D:1188:VAL:HG11	1.76	0.66
2:C:922:PHE:CE2	2:C:964:LYS:HB2	2.30	0.66
1:A:64:GLU:OE2	2:C:830:LYS:NZ	2.26	0.65
2:C:214:TYR:O	2:C:218:VAL:HG23	1.97	0.65
6:G:18:DT:C2	8:I:0:A:N1	2.64	0.65
1:A:106:PRO:HG3	1:A:134:GLU:HG2	1.77	0.65
5:F:370:LYS:HG2	5:F:375:LEU:HD23	1.77	0.65
3:D:832:ARG:HD2	3:D:833:GLU:H	1.62	0.65
5:F:400:ILE:HA	5:F:403:LYS:HG2	1.78	0.65
2:C:580:MET:HB3	2:C:584:GLU:CD	2.21	0.65
3:D:65:ARG:CG	3:D:66:GLN:H	2.10	0.65
3:D:1242:HIS:CD2	11:D:2005:UTP:O3B	2.50	0.65
5:F:274:THR:OG1	5:F:295:MET:HE3	1.97	0.65
1:B:112:ARG:NH1	1:B:126:ASP:OD1	2.30	0.65
2:C:627:ARG:CZ	2:C:640:ARG:HG3	2.26	0.65
3:D:266:GLU:OE2	3:D:315:ARG:HG3	1.97	0.64
3:D:1140:ILE:CG2	3:D:1144:LEU:HD12	2.27	0.64
3:D:237:LYS:N	3:D:240:GLU:OE1	2.29	0.64
2:C:1:MET:HB2	2:C:898:GLY:O	1.97	0.64
2:C:757:GLY:HA2	2:C:789:SER:OG	1.97	0.64
1:B:18:ARG:O	1:B:207:PRO:HD3	1.97	0.64
2:C:302:VAL:O	2:C:305:PRO:HD2	1.98	0.64
5:F:168:LYS:CG	5:F:172:ARG:HH21	2.09	0.64
2:C:480:THR:HG22	2:C:482:GLU:H	1.62	0.64
2:C:1070:ILE:CG2	3:D:655:PRO:HB2	2.28	0.64
1:B:65:PHE:CD2	3:D:809:PRO:HB2	2.32	0.64
2:C:740:GLU:HB3	2:C:805:ARG:HH12	1.63	0.63
3:D:236:TYR:CE1	3:D:242:LEU:HD12	2.33	0.63
3:D:65:ARG:HG2	3:D:66:GLN:H	1.63	0.63
4:E:70:THR:OG1	4:E:72:ARG:HG3	1.96	0.63
2:C:170:PRO:HD2	2:C:267:TYR:HE1	1.63	0.63
2:C:182:VAL:HG11	2:C:221:LEU:HG	1.81	0.63
5:F:329:TYR:CE2	5:F:333:ILE:HD11	2.34	0.63
1:B:44:LEU:HD13	1:B:199:ILE:HD13	1.79	0.63
1:B:124:ASN:ND2	1:B:127:LEU:HD22	2.13	0.63
2:C:704:HIS:CD2	2:C:831:ARG:HD2	2.33	0.63
2:C:911:GLU:OE1	3:D:1062:ARG:NH1	2.32	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1493:LYS:HA	3:D:1493:LYS:CE	2.23	0.63
2:C:709:GLU:OE2	2:C:824:ARG:NH1	2.32	0.63
2:C:769:PRO:CG	3:D:65:ARG:NH2	2.62	0.63
6:G:11:DG:C2'	6:G:12:DG:H5''	2.28	0.63
1:B:101:LEU:HB2	1:B:114:PHE:CD1	2.33	0.63
2:C:722:ILE:HD12	2:C:821:GLU:CD	2.24	0.63
5:F:274:THR:OG1	5:F:295:MET:CE	2.47	0.63
2:C:524:VAL:CG1	2:C:528:GLU:HB2	2.28	0.62
5:F:362:SER:OG	5:F:365:GLU:HG2	1.99	0.62
3:D:658:LEU:HA	3:D:661:MET:HE3	1.81	0.62
5:F:375:LEU:HD12	5:F:375:LEU:O	1.98	0.62
3:D:1156:LEU:HD23	3:D:1182:GLU:OE2	1.99	0.62
3:D:1492:LEU:HD13	3:D:1493:LYS:HZ1	1.64	0.62
3:D:1087:ARG:CB	3:D:1237:THR:CG2	2.78	0.62
3:D:1241:PHE:O	3:D:1243:THR:N	2.26	0.62
2:C:627:ARG:NH2	2:C:640:ARG:HG3	2.13	0.62
3:D:235:ALA:HB2	3:D:320:ALA:O	1.99	0.62
3:D:242:LEU:HB3	3:D:311:LEU:HD12	1.81	0.62
3:D:14:SER:HB3	3:D:511:TRP:CZ2	2.35	0.62
3:D:1258:ARG:HH21	3:D:1351:GLU:HG3	1.65	0.62
1:B:101:LEU:HD21	1:B:109:VAL:HG11	1.79	0.62
1:B:138:LEU:CD2	1:B:140:MET:HE3	2.27	0.62
3:D:1044:LEU:H	3:D:1044:LEU:HD12	1.65	0.62
3:D:34:TYR:CZ	3:D:35:ARG:HG3	2.34	0.61
3:D:223:LEU:HD11	3:D:288:MET:HE1	1.80	0.61
3:D:322:VAL:HG22	3:D:335:LEU:HD23	1.78	0.61
1:A:70:GLY:N	2:C:607:ASP:OD1	2.28	0.61
1:A:226:SER:O	1:A:228:PRO:HD3	1.99	0.61
2:C:764:GLU:O	2:C:766:GLU:N	2.28	0.61
3:D:65:ARG:HG2	3:D:66:GLN:N	2.16	0.61
4:E:50:THR:HG22	4:E:51:LEU:H	1.65	0.61
3:D:1493:LYS:HZ1	3:D:1496:GLU:HG3	1.64	0.61
3:D:405:ASP:N	3:D:423:ASP:OD1	2.25	0.61
5:F:193:ARG:NH1	7:H:7:DC:H5	1.99	0.61
5:F:265:VAL:O	5:F:269:ASN:ND2	2.30	0.61
6:G:17:DT:O2	8:I:1:A:C2	2.53	0.61
2:C:930:LYS:HE3	2:C:935:GLY:HA2	1.83	0.61
5:F:168:LYS:O	5:F:172:ARG:HG2	2.00	0.61
3:D:1108:ARG:HD2	3:D:1198:TYR:O	2.01	0.61
1:B:143:ARG:HG2	1:B:143:ARG:NH1	2.16	0.61
2:C:280:LYS:HE3	2:C:309:TYR:CZ	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1058:ASP:OD2	3:D:621:LYS:HE2	2.00	0.60
3:D:875:THR:HG21	3:D:902:LEU:HD21	1.82	0.60
5:F:354:LEU:HD23	5:F:418:LEU:HD21	1.83	0.60
3:D:1277:ILE:HG22	3:D:1278:ASP:H	1.65	0.60
5:F:274:THR:HG21	5:F:295:MET:HE3	1.83	0.60
3:D:284:LEU:CD2	3:D:288:MET:HE2	2.29	0.60
2:C:144:PRO:HG2	2:C:165:LEU:CD2	2.31	0.60
2:C:196:LEU:HD22	2:C:196:LEU:O	2.02	0.60
3:D:1112:CYS:SG	3:D:1114:THR:HG22	2.42	0.60
3:D:218:LYS:HG2	3:D:338:GLU:CG	2.21	0.60
3:D:1197:ARG:HG2	3:D:1198:TYR:CD2	2.37	0.60
3:D:236:TYR:CZ	3:D:242:LEU:HD12	2.37	0.60
3:D:17:LYS:HA	3:D:20:SER:OG	2.02	0.60
3:D:258:VAL:HG13	3:D:273:ARG:HG3	1.83	0.60
2:C:926:PHE:HE1	2:C:929:ARG:HH11	1.50	0.59
3:D:741:ASP:OD1	3:D:743:ASP:OD1	2.20	0.59
3:D:834:THR:OG1	3:D:838:ARG:HD2	2.02	0.59
1:A:97:VAL:HG13	1:A:98:THR:H	1.65	0.59
1:B:138:LEU:HD11	1:B:140:MET:CE	2.32	0.59
3:D:514:LEU:HD13	3:D:517:VAL:HG22	1.83	0.59
5:F:104:ARG:HG2	5:F:108:GLU:OE2	2.01	0.59
3:D:44:LEU:HB3	3:D:525:ARG:HH21	1.68	0.59
3:D:135:LEU:CD2	3:D:463:GLN:HG2	2.33	0.59
2:C:683:ASN:HB3	2:C:872:ASN:HD22	1.65	0.59
3:D:860:LEU:O	3:D:876:SER:HB2	2.02	0.59
3:D:1271:LYS:HG3	3:D:1272:ALA:O	2.01	0.59
3:D:1047:LYS:HB3	3:D:1048:PRO:CD	2.33	0.59
1:A:23:PHE:CD1	1:A:211:LEU:HD23	2.38	0.59
2:C:194:VAL:HG22	2:C:221:LEU:HD23	1.85	0.59
3:D:1044:LEU:HD12	3:D:1044:LEU:N	2.17	0.59
3:D:1347:TYR:O	3:D:1351:GLU:HB2	2.03	0.59
6:G:18:DT:N3	8:I:0:A:N6	2.07	0.59
2:C:134:ARG:NH1	2:C:392:SER:O	2.36	0.59
3:D:478:LEU:O	3:D:481:MET:HB2	2.03	0.59
3:D:809:PRO:HG3	3:D:829:VAL:HG11	1.83	0.59
3:D:1495:ILE:HG12	4:E:88:GLU:CD	2.28	0.59
5:F:101:GLU:HG2	5:F:105:LYS:HE2	1.84	0.59
2:C:32:ALA:HB2	2:C:73:LEU:CD1	2.30	0.58
2:C:394:PHE:CE2	2:C:632:ASN:CG	2.80	0.58
3:D:373:PRO:HA	3:D:376:GLU:OE2	2.04	0.58
3:D:963:TYR:CD2	3:D:1002:LYS:HD3	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:64:LEU:HD22	2:C:100:LEU:HD11	1.84	0.58
2:C:1059:ASP:OD1	2:C:1062:GLY:N	2.25	0.58
3:D:1042:ARG:HB3	3:D:1057:VAL:HB	1.85	0.58
2:C:212:GLY:CA	2:C:218:VAL:HG21	2.32	0.58
3:D:704:ARG:HB2	3:D:745:MET:HE2	1.84	0.58
3:D:1047:LYS:HB3	3:D:1048:PRO:HD2	1.83	0.58
3:D:823:LEU:HD11	3:D:837:GLY:HA2	1.85	0.58
6:G:18:DT:H2''	6:G:19:DG:O4'	2.04	0.58
2:C:376:ARG:HB2	2:C:377:PRO:HD3	1.85	0.58
3:D:314:PRO:HD2	3:D:317:VAL:HG13	1.86	0.58
3:D:455:ARG:HH12	3:D:463:GLN:HE21	1.52	0.58
5:F:329:TYR:HE2	5:F:333:ILE:HD11	1.69	0.58
2:C:776:SER:OG	5:F:373:LYS:NZ	2.36	0.58
3:D:64:LYS:HB3	5:F:378:GLY:HA3	1.85	0.58
2:C:32:ALA:CB	2:C:73:LEU:HD12	2.31	0.58
3:D:1208:ASP:OD1	3:D:1208:ASP:C	2.47	0.58
1:A:39:PRO:HG3	1:B:39:PRO:HG3	1.85	0.57
1:A:97:VAL:CG1	1:A:98:THR:N	2.67	0.57
6:G:18:DT:H2'	6:G:19:DG:H8	1.69	0.57
1:A:39:PRO:HG3	1:B:39:PRO:CG	2.33	0.57
2:C:480:THR:HG22	2:C:481:ASP:N	2.18	0.57
2:C:764:GLU:C	2:C:766:GLU:N	2.62	0.57
3:D:693:GLU:HB3	4:E:48:MET:HE1	1.87	0.57
5:F:96:LEU:O	5:F:100:VAL:HG23	2.04	0.57
5:F:82:ARG:HH21	7:H:9:DA:H61	1.52	0.57
2:C:638:ASP:OD1	2:C:640:ARG:HD2	2.04	0.57
2:C:740:GLU:HB3	2:C:805:ARG:NH1	2.19	0.57
2:C:143:SER:O	2:C:147:TYR:OH	2.18	0.57
2:C:150:PRO:HD3	2:C:322:VAL:HG11	1.86	0.57
2:C:999:HIS:HB3	2:C:1004:LYS:NZ	2.20	0.57
3:D:689:ASP:CG	4:E:51:LEU:HD11	2.30	0.57
7:H:17:DC:H2''	7:H:18:DG:H5'	1.86	0.57
1:A:137:ARG:HH11	1:A:137:ARG:HG2	1.70	0.56
2:C:359:MET:C	2:C:360:LEU:HD12	2.30	0.56
3:D:1154:GLU:HG2	3:D:1159:ARG:CA	2.31	0.56
2:C:261:ILE:HG21	2:C:291:ALA:HB3	1.88	0.56
2:C:429:ASP:OD1	3:D:1048:PRO:HB3	2.05	0.56
2:C:858:MET:HG2	2:C:867:VAL:O	2.05	0.56
3:D:355:VAL:HG11	3:D:385:VAL:HG21	1.88	0.56
2:C:396:ASP:HA	2:C:633:GLN:NE2	2.20	0.56
2:C:766:GLU:CB	2:C:767:PRO:CD	2.83	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:144:PRO:HG2	2:C:165:LEU:HD23	1.87	0.56
2:C:240:THR:HG22	2:C:248:PRO:HG3	1.88	0.56
1:A:64:GLU:HG2	1:A:76:VAL:HG22	1.87	0.56
2:C:987:ILE:HD11	3:D:946:GLY:HA2	1.88	0.56
3:D:1284:GLU:HG2	3:D:1291:SER:HB3	1.88	0.56
3:D:664:LYS:HB2	3:D:666:ILE:HG13	1.88	0.56
2:C:614:ARG:HH21	2:C:614:ARG:HG3	1.70	0.56
3:D:181:ASP:HB2	3:D:205:TYR:CG	2.40	0.56
3:D:828:LYS:HG2	3:D:833:GLU:HB3	1.88	0.56
5:F:101:GLU:OE1	5:F:105:LYS:HE3	2.06	0.56
1:B:83:LYS:HE2	1:B:168:ASP:HB2	1.86	0.56
2:C:182:VAL:HG21	2:C:193:LEU:HD23	1.87	0.56
5:F:274:THR:CB	5:F:295:MET:HE3	2.36	0.56
3:D:1492:LEU:HD13	3:D:1493:LYS:NZ	2.20	0.56
3:D:131:LYS:O	3:D:456:MET:HG2	2.05	0.55
3:D:983:LEU:HA	3:D:987:GLU:OE2	2.07	0.55
2:C:150:PRO:HA	2:C:158:TYR:HD1	1.72	0.55
1:A:198:ARG:HD3	2:C:934:PHE:CZ	2.42	0.55
2:C:1001:VAL:HG13	3:D:630:VAL:HB	1.88	0.55
3:D:996:TRP:HB3	3:D:1044:LEU:HD11	1.88	0.55
3:D:1444:THR:O	3:D:1448:THR:HG23	2.05	0.55
3:D:564:GLU:OE2	5:F:140:ARG:NH2	2.40	0.55
3:D:907:GLU:OE1	3:D:909:ASN:N	2.38	0.55
2:C:182:VAL:CG1	2:C:221:LEU:HG	2.37	0.55
2:C:214:TYR:HB3	2:C:217:LEU:HD12	1.89	0.55
3:D:1036:ARG:HH22	3:D:1246:VAL:HG11	1.71	0.55
3:D:984:THR:OG1	3:D:987:GLU:HG3	2.07	0.55
3:D:1386:ASP:OD2	3:D:1413:THR:HG22	2.07	0.55
5:F:122:LEU:HB2	5:F:127:ILE:HD11	1.89	0.55
6:G:20:DT:H2'	6:G:20:DT:O2	2.06	0.55
2:C:1043:TYR:CD1	3:D:763:MET:HG2	2.42	0.54
3:D:65:ARG:CG	3:D:66:GLN:N	2.69	0.54
3:D:1336:LEU:HB2	3:D:1344:VAL:HG21	1.88	0.54
2:C:370:ALA:O	5:F:280:GLN:NE2	2.29	0.54
5:F:271:LEU:HD22	5:F:307:THR:HG21	1.89	0.54
5:F:287:THR:HG22	5:F:288:TYR:H	1.71	0.54
3:D:231:VAL:HG13	3:D:242:LEU:O	2.07	0.54
3:D:1339:LYS:HB3	3:D:1343:ALA:CB	2.37	0.54
3:D:1461:GLY:O	3:D:1465:ASN:ND2	2.40	0.54
3:D:1492:LEU:CD1	3:D:1493:LYS:NZ	2.70	0.54
2:C:577:PRO:HG2	2:C:580:MET:HG2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1059:ASP:OD1	2:C:1059:ASP:C	2.50	0.54
3:D:1046:GLN:HA	3:D:1052:THR:HA	1.89	0.54
1:A:206:THR:HB	1:A:209:GLU:HG3	1.89	0.54
2:C:774:LEU:HD22	5:F:350:LEU:HD11	1.90	0.54
1:B:80:LEU:HD13	3:D:867:ARG:HD3	1.90	0.54
2:C:880:MET:HB3	3:D:1061:PHE:CE2	2.42	0.54
3:D:1386:ASP:H	3:D:1413:THR:HG23	1.73	0.54
3:D:1252:ILE:HG23	3:D:1252:ILE:O	2.07	0.54
1:B:179:PHE:HB3	1:B:197:LEU:CD2	2.38	0.54
3:D:236:TYR:HA	3:D:240:GLU:OE1	2.08	0.54
3:D:1040:GLY:O	3:D:1060:SER:HB3	2.07	0.54
4:E:12:MET:HE1	4:E:69:LEU:HA	1.90	0.54
2:C:211:LEU:HD23	2:C:311:PHE:CD2	2.43	0.53
1:B:26:GLU:HB3	1:B:194:LYS:HG3	1.91	0.53
2:C:1058:ASP:OD1	2:C:1084:SER:HB2	2.08	0.53
3:D:231:VAL:O	3:D:236:TYR:OH	2.25	0.53
3:D:520:LEU:O	3:D:525:ARG:NH1	2.39	0.53
3:D:963:TYR:HE2	3:D:1002:LYS:HD3	1.67	0.53
5:F:345:ALA:O	5:F:349:LEU:HG	2.08	0.53
2:C:170:PRO:HD2	2:C:267:TYR:CD1	2.43	0.53
2:C:261:ILE:HG22	2:C:262:ALA:N	2.22	0.53
2:C:394:PHE:CE2	2:C:632:ASN:CB	2.90	0.53
3:D:59:ALA:HB2	3:D:78:VAL:HG21	1.90	0.53
3:D:1154:GLU:CG	3:D:1159:ARG:HA	2.34	0.53
1:B:102:LYS:HG2	1:B:139:ASN:OD1	2.08	0.53
2:C:614:ARG:HG3	2:C:614:ARG:NH2	2.24	0.53
3:D:111:LYS:HG3	3:D:1452:ILE:HD11	1.89	0.53
1:A:106:PRO:CG	1:A:134:GLU:HG2	2.38	0.53
1:B:73:GLU:OE1	1:B:73:GLU:N	2.30	0.53
3:D:1084:THR:CG2	3:D:1238:MET:HA	2.38	0.53
1:A:100:LEU:CD2	1:A:141:GLU:HG2	2.39	0.53
2:C:805:ARG:HG3	2:C:823:VAL:HG22	1.90	0.53
1:A:196:THR:HG21	2:C:934:PHE:CE2	2.44	0.53
2:C:154:ARG:HH21	2:C:157:ARG:HG3	1.74	0.53
2:C:617:ASP:OD1	2:C:617:ASP:N	2.41	0.53
3:D:1154:GLU:CD	3:D:1159:ARG:HB2	2.33	0.53
3:D:1276:GLU:O	3:D:1277:ILE:HD13	2.08	0.53
5:F:368:VAL:HG21	5:F:400:ILE:HD11	1.90	0.53
2:C:150:PRO:HD3	2:C:322:VAL:CG1	2.39	0.53
3:D:664:LYS:C	3:D:666:ILE:H	2.17	0.52
3:D:1191:PRO:HD2	3:D:1369:GLU:OE1	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1276:GLU:C	3:D:1277:ILE:HD13	2.34	0.52
2:C:329:GLY:O	2:C:466:PHE:CE2	2.62	0.52
2:C:374:ASN:CG	5:F:276:ARG:HD2	2.33	0.52
2:C:1100:GLN:HG3	3:D:9:ARG:HH21	1.75	0.52
2:C:258:TYR:CD2	2:C:262:ALA:HB3	2.44	0.52
2:C:132:ALA:HB1	2:C:394:PHE:CE1	2.41	0.52
3:D:1386:ASP:HB3	3:D:1412:LYS:HD2	1.91	0.52
2:C:593:ALA:HB1	2:C:659:PRO:HD2	1.92	0.52
1:B:179:PHE:HB3	1:B:197:LEU:HD23	1.91	0.52
3:D:15:PRO:O	3:D:19:ARG:HG3	2.09	0.52
3:D:218:LYS:CG	3:D:338:GLU:HG2	2.23	0.52
2:C:84:ARG:HD3	2:C:629:TYR:OH	2.09	0.52
3:D:114:THR:CG2	3:D:498:VAL:HG21	2.39	0.52
3:D:892:ASP:OD1	3:D:894:LYS:HD2	2.08	0.52
2:C:462:ASP:OD2	2:C:468:ARG:NH1	2.40	0.52
1:A:112:ARG:HG2	1:A:125:PRO:O	2.09	0.52
2:C:359:MET:O	2:C:360:LEU:HD12	2.10	0.52
6:G:18:DT:C2'	6:G:19:DG:O4'	2.57	0.52
3:D:495:ARG:O	3:D:499:VAL:HG23	2.10	0.52
1:A:44:LEU:O	1:A:174:VAL:HG21	2.10	0.51
2:C:260:LEU:O	2:C:261:ILE:HD12	2.10	0.51
2:C:575:GLN:HG3	2:C:670:GLN:HA	1.92	0.51
3:D:135:LEU:HD23	3:D:463:GLN:HG2	1.91	0.51
3:D:317:VAL:HG23	3:D:339:TRP:HB3	1.91	0.51
3:D:899:LEU:HD22	3:D:917:GLN:HB3	1.93	0.51
3:D:1281:VAL:HG23	3:D:1317:ASP:O	2.10	0.51
2:C:418:LEU:H	2:C:418:LEU:HD12	1.75	0.51
1:B:90:LEU:O	1:B:92:PRO:HD3	2.10	0.51
3:D:314:PRO:HB2	3:D:317:VAL:CG1	2.40	0.51
2:C:587:VAL:O	2:C:591:SER:HB3	2.09	0.51
3:D:134:VAL:HG23	3:D:149:LYS:HA	1.93	0.51
3:D:1137:ARG:O	3:D:1141:GLU:HG3	2.10	0.51
6:G:17:DT:N3	8:I:1:A:C6	2.61	0.51
6:G:18:DT:H2''	6:G:19:DG:C5'	2.39	0.51
1:A:23:PHE:N	1:A:23:PHE:CD2	2.77	0.51
5:F:260:ILE:HG22	5:F:265:VAL:HG23	1.92	0.51
1:B:188:GLN:HG2	1:B:189:ARG:N	2.26	0.51
3:D:945:SER:OG	3:D:947:ILE:HG12	2.11	0.51
3:D:1263:PHE:HA	3:D:1375:MET:CE	2.38	0.51
6:G:9:DC:H2''	6:G:10:DC:OP2	2.11	0.51
1:B:47:SER:HB3	1:B:217:ILE:HD13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:269:LEU:O	2:C:288:ARG:HD2	2.11	0.51
3:D:817:GLU:OE1	3:D:817:GLU:HA	2.06	0.51
2:C:136:ILE:CB	2:C:336:VAL:HG13	2.40	0.51
3:D:204:LEU:HD22	3:D:441:ARG:NH1	2.25	0.51
5:F:222:ARG:HA	5:F:222:ARG:CZ	2.40	0.51
6:G:12:DG:H8	6:G:12:DG:C5'	2.23	0.51
5:F:393:THR:HG22	5:F:395:GLU:N	2.24	0.51
2:C:758:ARG:HH21	2:C:788:THR:HB	1.75	0.50
2:C:760:SER:HB2	2:C:786:LYS:HE3	1.93	0.50
2:C:765:SER:O	2:C:766:GLU:O	2.29	0.50
1:A:11:PHE:CD1	1:B:225:PHE:HA	2.47	0.50
1:A:191:ASP:OD1	2:C:938:LYS:NZ	2.29	0.50
2:C:230:ARG:HG3	2:C:233:GLU:HG3	1.93	0.50
2:C:499:ALA:HB2	2:C:533:ASP:HB2	1.93	0.50
2:C:563:ASN:O	2:C:566:THR:HB	2.12	0.50
3:D:84:ILE:CD1	3:D:87:ARG:HD2	2.42	0.50
1:A:23:PHE:CD1	1:A:211:LEU:CD2	2.95	0.50
1:A:57:TYR:CG	1:A:161:ARG:HD2	2.46	0.50
3:D:693:GLU:CB	4:E:48:MET:HE1	2.41	0.50
3:D:520:LEU:HD11	3:D:524:LEU:HD12	1.94	0.50
3:D:796:ARG:NH2	3:D:859:ASP:OD2	2.34	0.50
3:D:1493:LYS:HE3	3:D:1496:GLU:HB2	1.92	0.50
1:B:80:LEU:HD22	3:D:844:ALA:CA	2.40	0.50
4:E:46:PRO:HB2	4:E:57:ASP:CB	2.41	0.50
2:C:340:MET:CE	2:C:382:ILE:HD13	2.40	0.50
3:D:142:LEU:HB2	3:D:161:LEU:HD11	1.93	0.50
3:D:1313:VAL:HG11	3:D:1319:VAL:HG11	1.92	0.50
3:D:236:TYR:HB3	3:D:313:MET:HG3	1.94	0.50
1:B:190:THR:HG22	1:B:190:THR:O	2.12	0.50
2:C:468:ARG:HA	2:C:486:MET:O	2.10	0.50
2:C:722:ILE:HD12	2:C:821:GLU:HG3	1.93	0.50
3:D:526:PRO:HB2	3:D:528:VAL:HG13	1.93	0.50
1:B:94:LEU:HD21	1:B:97:VAL:CG2	2.42	0.50
3:D:780:LYS:HE3	3:D:912:LYS:HD3	1.94	0.50
2:C:361:MET:O	2:C:362:GLY:O	2.29	0.49
2:C:591:SER:O	2:C:592:LEU:HB2	2.12	0.49
2:C:888:THR:HG22	2:C:990:GLY:HA3	1.93	0.49
2:C:911:GLU:CD	3:D:1062:ARG:HH11	2.19	0.49
3:D:805:GLU:HG3	3:D:828:LYS:HB2	1.94	0.49
1:B:108:GLU:CG	1:B:131:THR:HG22	2.34	0.49
3:D:215:TYR:O	3:D:340:THR:HA	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:750:PRO:O	3:D:756:GLN:NE2	2.45	0.49
2:C:204:GLN:HB2	2:C:227:PHE:CZ	2.47	0.49
2:C:235:LEU:HD21	2:C:254:VAL:HG22	1.93	0.49
2:C:482:GLU:HG3	2:C:482:GLU:O	2.12	0.49
2:C:693:GLU:HG2	2:C:855:VAL:HB	1.95	0.49
2:C:861:LEU:HB3	2:C:862:PRO:HD2	1.94	0.49
3:D:351:MET:HG2	3:D:370:ALA:HB2	1.94	0.49
5:F:274:THR:CG2	5:F:295:MET:HE3	2.42	0.49
1:A:39:PRO:CG	1:B:39:PRO:HG3	2.42	0.49
2:C:216:GLU:HA	2:C:216:GLU:OE1	2.12	0.49
2:C:343:GLN:HA	2:C:343:GLN:NE2	2.26	0.49
3:D:231:VAL:O	3:D:231:VAL:HG22	2.12	0.49
3:D:796:ARG:HH22	3:D:859:ASP:CG	2.20	0.49
2:C:150:PRO:HG3	2:C:322:VAL:HG21	1.94	0.49
2:C:405:ARG:O	2:C:405:ARG:HG2	2.12	0.49
2:C:598:GLU:O	2:C:651:LYS:HG3	2.12	0.49
1:A:89:PHE:HD2	1:A:94:LEU:O	1.95	0.49
1:A:215:VAL:HG13	1:B:222:LEU:HD22	1.95	0.49
2:C:603:VAL:HG23	2:C:647:GLN:O	2.13	0.49
2:C:1056:LYS:HG2	3:D:625:TYR:HD1	1.77	0.49
3:D:215:TYR:HE1	3:D:381:ALA:H	1.58	0.49
3:D:1208:ASP:OD2	3:D:1211:MET:HE2	2.13	0.49
3:D:1364:HIS:ND1	3:D:1366:LYS:CB	2.75	0.49
1:A:96:THR:HB	1:A:145:ASP:OD1	2.12	0.49
2:C:77:PRO:HB2	2:C:78:PHE:CD1	2.47	0.49
2:C:396:ASP:HA	2:C:633:GLN:HE22	1.78	0.49
2:C:571:LEU:HD22	2:C:700:TYR:HA	1.93	0.49
3:D:270:LEU:HD23	3:D:270:LEU:C	2.36	0.49
3:D:1283:ILE:CG1	3:D:1315:ASP:CG	2.84	0.49
3:D:1353:GLN:HA	3:D:1353:GLN:OE1	2.12	0.49
3:D:65:ARG:HH11	3:D:65:ARG:CB	2.19	0.49
3:D:1141:GLU:CD	3:D:1168:MET:HE1	2.36	0.49
3:D:1377:LYS:HE3	3:D:1378:TYR:CZ	2.47	0.49
4:E:54:LEU:C	4:E:55:PHE:HD2	2.20	0.49
7:H:7:DC:H6	7:H:7:DC:H5'	1.78	0.49
2:C:55:GLU:O	2:C:359:MET:HE1	2.13	0.49
2:C:679:PHE:H	2:C:683:ASN:HD22	1.59	0.49
2:C:1053:LEU:HA	3:D:621:LYS:HD2	1.95	0.49
2:C:539:VAL:O	2:C:539:VAL:HG12	2.12	0.48
2:C:679:PHE:N	2:C:683:ASN:HD22	2.10	0.48
3:D:1263:PHE:HD2	3:D:1375:MET:HE3	1.75	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:278:LEU:HD11	5:F:294:ALA:CB	2.43	0.48
3:D:103:TRP:CZ2	3:D:1444:THR:HG22	2.49	0.48
5:F:193:ARG:NH1	7:H:7:DC:C5	2.79	0.48
3:D:1386:ASP:N	3:D:1413:THR:HG23	2.28	0.48
5:F:238:TYR:OH	7:H:1:DT:H73	2.14	0.48
5:F:279:GLN:O	5:F:279:GLN:HG2	2.08	0.48
5:F:353:GLU:HB3	5:F:418:LEU:HD11	1.95	0.48
1:B:111:ALA:HB3	1:B:125:PRO:HA	1.95	0.48
2:C:304:LEU:HB3	2:C:305:PRO:HD3	1.95	0.48
3:D:465:LEU:HD12	3:D:513:ILE:HD13	1.95	0.48
5:F:323:ASP:CB	6:G:19:DG:H1	2.26	0.48
2:C:717:LEU:CD2	2:C:763:GLY:HA2	2.44	0.48
2:C:774:LEU:HD21	5:F:421:PHE:HB2	1.96	0.48
2:C:952:LEU:HA	2:C:952:LEU:HD23	1.64	0.48
2:C:999:HIS:HB3	2:C:1004:LYS:HZ2	1.78	0.48
3:D:103:TRP:HB3	3:D:1448:THR:HG21	1.95	0.48
5:F:400:ILE:HA	5:F:403:LYS:CG	2.41	0.48
2:C:394:PHE:CZ	2:C:632:ASN:CG	2.91	0.48
2:C:713:ARG:CZ	3:D:533:GLY:HA3	2.44	0.48
3:D:654:LYS:HB3	3:D:655:PRO:HD3	1.95	0.48
5:F:383:LEU:HD12	5:F:394:ARG:NH2	2.28	0.48
1:A:57:TYR:HE1	1:A:163:ASN:HB2	1.78	0.48
3:D:1410:GLU:O	3:D:1410:GLU:HG2	2.14	0.48
5:F:154:LYS:O	5:F:158:GLU:HG3	2.13	0.48
2:C:460:ARG:CD	2:C:485:TYR:CZ	2.94	0.48
3:D:792:ILE:HD13	3:D:941:PHE:CE1	2.49	0.48
3:D:1339:LYS:HB3	3:D:1343:ALA:HB3	1.96	0.48
1:A:198:ARG:HD3	2:C:934:PHE:CE1	2.49	0.48
2:C:267:TYR:CE2	2:C:290:LEU:HG	2.49	0.48
3:D:147:VAL:HG21	3:D:161:LEU:HD21	1.96	0.48
3:D:1047:LYS:HG2	3:D:1053:PHE:CZ	2.49	0.48
1:B:206:THR:CG2	1:B:209:GLU:H	2.16	0.48
2:C:787:ASP:OD2	2:C:791:ARG:NH2	2.47	0.48
3:D:792:ILE:HD13	3:D:941:PHE:CD1	2.49	0.48
3:D:1486:VAL:HG22	4:E:22:VAL:HG13	1.93	0.48
1:A:128:HIS:O	1:A:129:ILE:HD13	2.13	0.47
1:B:20:TYR:OH	1:B:22:GLU:HG2	2.13	0.47
2:C:260:LEU:C	2:C:261:ILE:HD12	2.39	0.47
3:D:1068:LEU:O	3:D:1072:ILE:HG12	2.14	0.47
1:B:49:PRO:CA	1:B:148:VAL:HG12	2.38	0.47
2:C:486:MET:HE1	2:C:531:PHE:CE1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1101:THR:OG1	2:C:1109:VAL:O	2.33	0.47
3:D:1495:ILE:HD11	4:E:88:GLU:HB2	1.97	0.47
1:B:162:ILE:O	1:B:163:ASN:HB2	2.13	0.47
2:C:249:LYS:HB3	2:C:252:LYS:HB2	1.95	0.47
2:C:775:ARG:HH21	5:F:422:LEU:HD23	1.80	0.47
2:C:862:PRO:HA	2:C:975:TYR:CE2	2.49	0.47
3:D:181:ASP:HB2	3:D:205:TYR:CD1	2.49	0.47
3:D:1258:ARG:NH2	3:D:1351:GLU:HG3	2.28	0.47
2:C:26:TYR:OH	2:C:119:PRO:O	2.14	0.47
2:C:168:ARG:O	2:C:267:TYR:HA	2.13	0.47
2:C:243:ARG:HD2	2:C:256:TYR:CE1	2.49	0.47
3:D:1051:GLU:O	3:D:1051:GLU:HG2	2.11	0.47
3:D:1240:THR:CG2	3:D:1251:ASP:OD2	2.63	0.47
3:D:1249:ALA:O	3:D:1250:ALA:HB3	2.15	0.47
3:D:1342:GLU:CD	3:D:1342:GLU:H	2.22	0.47
3:D:1384:PRO:HB3	3:D:1389:LEU:O	2.13	0.47
5:F:82:ARG:HB2	7:H:8:DG:O6	2.14	0.47
5:F:358:LEU:HD13	5:F:370:LYS:HE3	1.96	0.47
2:C:480:THR:HG22	2:C:481:ASP:H	1.78	0.47
3:D:126:VAL:O	3:D:457:GLY:N	2.42	0.47
8:I:0:A:HO2'	8:I:1:A:P	2.31	0.47
1:A:47:SER:O	1:A:49:PRO:HD3	2.14	0.47
2:C:13:ILE:HD13	2:C:483:VAL:HG21	1.97	0.47
5:F:130:VAL:HG21	5:F:159:ILE:CG2	2.41	0.47
5:F:277:GLN:O	5:F:281:GLU:HG2	2.14	0.47
1:A:24:VAL:HG22	1:A:196:THR:HG23	1.96	0.47
1:A:206:THR:HG22	1:A:208:LEU:N	2.29	0.47
2:C:181:VAL:HA	2:C:220:GLY:O	2.15	0.47
2:C:214:TYR:CB	2:C:217:LEU:HD12	2.45	0.47
2:C:1067:TYR:CE2	2:C:1071:ILE:HD13	2.50	0.47
3:D:179:VAL:HB	3:D:183:GLU:CD	2.39	0.47
3:D:407:VAL:HG12	5:F:172:ARG:HH12	1.78	0.47
3:D:776:GLU:OE2	3:D:1362:LYS:HD3	2.14	0.47
3:D:876:SER:OG	3:D:879:ARG:HG3	2.15	0.47
3:D:1047:LYS:CG	3:D:1053:PHE:CE1	2.96	0.47
2:C:711:GLU:OE2	2:C:816:LYS:HE3	2.14	0.47
5:F:101:GLU:OE1	5:F:105:LYS:CE	2.63	0.47
5:F:194:LEU:HB2	7:H:6:DT:O2	2.15	0.47
1:A:39:PRO:CG	1:B:39:PRO:CG	2.92	0.47
1:A:137:ARG:HG2	1:A:137:ARG:NH1	2.29	0.47
2:C:472:ARG:HD2	2:C:480:THR:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:976:ASP:OD1	2:C:978:ARG:HD3	2.14	0.47
3:D:236:TYR:HB3	3:D:313:MET:CG	2.44	0.47
3:D:422:ALA:H	3:D:427:VAL:HG13	1.80	0.47
3:D:1263:PHE:O	3:D:1375:MET:HE2	2.15	0.47
5:F:95:THR:HG22	5:F:96:LEU:N	2.30	0.47
8:I:2:U:C2'	8:I:3:U:H5'	2.45	0.47
1:A:38:ASN:HB3	1:A:39:PRO:HD3	1.96	0.46
3:D:871:LYS:NZ	3:D:897:TRP:HE1	2.13	0.46
3:D:1014:ASN:OD1	3:D:1014:ASN:N	2.48	0.46
2:C:874:LEU:HD23	3:D:1023:MET:SD	2.55	0.46
2:C:69:LEU:HB2	2:C:97:ARG:O	2.14	0.46
3:D:175:VAL:CG1	3:D:193:PRO:HD2	2.45	0.46
3:D:586:ARG:NH2	6:G:10:DC:H5''	2.30	0.46
3:D:895:VAL:HG11	3:D:922:LEU:HD21	1.96	0.46
3:D:959:GLU:O	3:D:963:TYR:HD1	1.99	0.46
3:D:983:LEU:HB3	3:D:987:GLU:HB2	1.98	0.46
3:D:1122:LEU:HD22	3:D:1140:ILE:HD13	1.97	0.46
3:D:1272:ALA:HA	3:D:1326:THR:HB	1.96	0.46
5:F:197:SER:HA	5:F:200:LYS:HE2	1.98	0.46
1:A:6:LEU:HD11	1:A:192:LEU:HD13	1.98	0.46
1:B:65:PHE:HD2	3:D:809:PRO:HB2	1.78	0.46
2:C:194:VAL:CG1	2:C:221:LEU:O	2.64	0.46
2:C:419:THR:HG22	2:C:420:ARG:H	1.80	0.46
2:C:103:LYS:HD2	2:C:104:ASP:H	1.80	0.46
2:C:114:PHE:CD2	5:F:283:GLY:HA2	2.50	0.46
3:D:586:ARG:O	3:D:586:ARG:NH1	2.48	0.46
3:D:643:GLY:HA3	3:D:727:GLN:HB2	1.97	0.46
3:D:1100:ASP:OD2	3:D:1463:LYS:NZ	2.35	0.46
2:C:99:GLN:O	2:C:99:GLN:HG3	2.15	0.46
2:C:428:ARG:NH2	2:C:447:ALA:O	2.48	0.46
2:C:466:PHE:N	2:C:466:PHE:CD1	2.84	0.46
6:G:18:DT:C2'	6:G:19:DG:H5'	2.40	0.46
2:C:182:VAL:CG2	2:C:193:LEU:HD23	2.45	0.46
2:C:194:VAL:HG11	2:C:221:LEU:O	2.16	0.46
1:A:185:ARG:NH2	1:A:187:GLY:O	2.49	0.46
1:A:220:GLU:O	1:A:223:THR:OG1	2.25	0.46
2:C:164:PRO:CA	2:C:269:LEU:HD12	2.43	0.46
3:D:1161:GLU:CD	3:D:1164:ARG:HB2	2.41	0.46
5:F:140:ARG:HE	5:F:140:ARG:HB2	1.52	0.46
2:C:136:ILE:HD13	2:C:392:SER:HA	1.97	0.46
2:C:290:LEU:O	2:C:301:GLU:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:322:VAL:CG1	2:C:323:ASP:N	2.79	0.46
2:C:769:PRO:HG3	3:D:65:ARG:CZ	2.46	0.46
3:D:618:LEU:HD12	3:D:618:LEU:HA	1.64	0.46
1:B:153:ALA:C	1:B:155:LYS:H	2.24	0.46
2:C:162:ILE:HG21	2:C:290:LEU:HD21	1.98	0.46
2:C:911:GLU:OE2	3:D:1062:ARG:HD3	2.16	0.46
3:D:4:GLU:OE2	3:D:6:ARG:HD3	2.16	0.46
3:D:156:GLU:CD	3:D:156:GLU:H	2.24	0.46
3:D:1263:PHE:CD2	3:D:1375:MET:CE	2.94	0.46
3:D:46:ASP:OD1	3:D:48:ARG:HB2	2.17	0.45
3:D:361:VAL:O	3:D:382:GLU:HA	2.16	0.45
1:B:79:ILE:HA	1:B:82:LEU:HD12	1.97	0.45
1:B:95:GLN:NE2	1:B:146:ARG:HD2	2.31	0.45
3:D:471:GLU:H	3:D:471:GLU:HG2	1.58	0.45
1:B:208:LEU:HD12	1:B:208:LEU:O	2.16	0.45
2:C:722:ILE:HD12	2:C:821:GLU:CG	2.47	0.45
2:C:838:LYS:HE3	3:D:741:ASP:O	2.17	0.45
2:C:882:LEU:HD23	2:C:882:LEU:HA	1.66	0.45
3:D:553:ARG:HD2	3:D:570:GLU:OE2	2.15	0.45
2:C:198:ARG:NH2	2:C:203:ASP:OD1	2.49	0.45
1:A:206:THR:HG22	1:A:209:GLU:H	1.81	0.45
3:D:103:TRP:HB3	3:D:1448:THR:CG2	2.46	0.45
3:D:455:ARG:HH12	3:D:463:GLN:NE2	2.14	0.45
1:B:95:GLN:O	1:B:96:THR:HB	2.17	0.45
2:C:797:GLY:O	2:C:829:GLN:NE2	2.49	0.45
3:D:135:LEU:O	3:D:149:LYS:HE3	2.17	0.45
3:D:486:ARG:HA	3:D:489:ARG:NH2	2.31	0.45
3:D:1248:GLY:O	3:D:1249:ALA:O	2.33	0.45
3:D:1261:GLU:OE1	3:D:1268:PRO:HA	2.16	0.45
3:D:1265:ALA:O	3:D:1333:HIS:HE1	2.00	0.45
1:A:22:GLU:C	1:A:23:PHE:CD2	2.95	0.45
1:B:141:GLU:OE1	1:B:161:ARG:NH2	2.50	0.45
2:C:184:MET:HB2	2:C:193:LEU:HD13	1.99	0.45
2:C:226:VAL:HA	2:C:229:MET:HE2	1.99	0.45
2:C:593:ALA:HB1	2:C:659:PRO:CD	2.47	0.45
3:D:134:VAL:CG2	3:D:151:GLN:H	2.30	0.45
3:D:972:LEU:HD12	3:D:972:LEU:HA	1.64	0.45
3:D:1092:GLY:HA3	6:G:14:DA:O4'	2.16	0.45
2:C:44:ILE:HD13	2:C:44:ILE:HA	1.61	0.45
2:C:206:THR:O	2:C:210:GLU:HB2	2.16	0.45
2:C:501:THR:HG21	2:C:513:VAL:HB	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1364:HIS:CE1	3:D:1366:LYS:HB2	2.50	0.45
1:B:78:ILE:HG21	1:B:140:MET:HE1	1.98	0.45
3:D:17:LYS:HE2	3:D:17:LYS:HB2	1.58	0.45
3:D:804:LEU:HB2	3:D:806:PHE:HE2	1.81	0.45
5:F:207:LEU:HD23	5:F:207:LEU:HA	1.68	0.45
1:A:107:LYS:HE3	1:A:107:LYS:HB2	1.75	0.45
1:B:71:VAL:HG22	1:B:132:LEU:HG	1.99	0.45
2:C:470:PRO:CD	2:C:485:TYR:CE2	2.97	0.45
3:D:1144:LEU:O	3:D:1147:ARG:HG3	2.17	0.45
3:D:1379:VAL:HG21	3:D:1400:VAL:HG11	1.99	0.45
1:B:13:VAL:HG13	1:B:23:PHE:CE1	2.52	0.44
1:B:59:GLU:HB2	1:B:139:ASN:HB3	2.00	0.44
2:C:293:PHE:HD1	2:C:298:PHE:CE2	2.35	0.44
3:D:1036:ARG:NH2	3:D:1042:ARG:O	2.44	0.44
5:F:127:ILE:O	5:F:131:VAL:HG23	2.17	0.44
1:A:32:PHE:HA	1:A:35:THR:HB	1.99	0.44
3:D:399:ARG:NH1	3:D:448:GLU:OE2	2.51	0.44
3:D:880:ILE:HD13	3:D:880:ILE:HA	1.85	0.44
3:D:269:PHE:CE2	3:D:283:PHE:HD1	2.35	0.44
3:D:1216:SER:OG	4:E:16:LYS:HG2	2.17	0.44
3:D:1492:LEU:CD1	3:D:1493:LYS:HZ1	2.28	0.44
5:F:114:LYS:O	5:F:117:SER:HB3	2.17	0.44
1:A:100:LEU:HD23	1:A:141:GLU:HG2	2.00	0.44
1:A:186:LEU:O	1:A:188:GLN:OE1	2.36	0.44
1:B:80:LEU:HB3	3:D:867:ARG:HH21	1.81	0.44
2:C:340:MET:HE2	2:C:382:ILE:CD1	2.45	0.44
2:C:607:ASP:HB3	2:C:610:ARG:H	1.82	0.44
3:D:851:LEU:HD11	3:D:855:HIS:HE1	1.82	0.44
3:D:1478:SER:O	3:D:1482:ARG:HB2	2.17	0.44
5:F:229:TYR:CZ	5:F:230:LYS:HD3	2.52	0.44
1:A:29:GLU:O	1:A:32:PHE:HB2	2.17	0.44
2:C:580:MET:SD	2:C:584:GLU:HG3	2.57	0.44
2:C:704:HIS:HE1	2:C:1000:MET:HE1	1.80	0.44
3:D:204:LEU:HD23	3:D:204:LEU:HA	1.85	0.44
3:D:632:VAL:O	3:D:727:GLN:HA	2.16	0.44
1:A:11:PHE:CE1	1:B:225:PHE:HA	2.53	0.44
2:C:715:THR:HG22	2:C:716:LYS:N	2.32	0.44
3:D:108:VAL:HA	3:D:109:PRO:C	2.43	0.44
3:D:174:GLY:HA2	3:D:389:GLU:OE2	2.18	0.44
5:F:82:ARG:NH2	7:H:9:DA:H61	2.15	0.44
2:C:100:LEU:HD23	2:C:367:LEU:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:12:LEU:HD23	3:D:12:LEU:HA	1.79	0.44
3:D:1232:PRO:O	3:D:1236:LEU:HG	2.18	0.44
3:D:1353:GLN:NE2	3:D:1365:ASP:OD2	2.38	0.44
5:F:361:LEU:HB2	5:F:366:ALA:HB2	1.99	0.44
6:G:19:DG:H2''	6:G:20:DT:O5'	2.18	0.44
2:C:462:ASP:HB3	2:C:468:ARG:HD2	2.00	0.44
3:D:1197:ARG:HB2	3:D:1398:TRP:CZ2	2.52	0.44
3:D:1241:PHE:C	3:D:1243:THR:H	2.21	0.44
3:D:1386:ASP:N	3:D:1413:THR:CG2	2.81	0.44
5:F:239:ALA:O	5:F:240:THR:C	2.61	0.44
1:B:85:LEU:HG	1:B:87:VAL:HG23	1.99	0.44
2:C:374:ASN:OD1	2:C:376:ARG:HG2	2.18	0.44
2:C:659:PRO:O	2:C:660:ALA:C	2.61	0.44
3:D:191:LEU:HB2	3:D:195:VAL:HG12	2.00	0.44
3:D:473:LEU:HD21	3:D:495:ARG:HH21	1.81	0.44
4:E:14:ASP:OD2	4:E:18:ARG:NH1	2.45	0.44
1:A:57:TYR:CE1	1:A:163:ASN:HB2	2.53	0.43
3:D:468:LEU:HD23	3:D:468:LEU:HA	1.77	0.43
2:C:269:LEU:HD12	2:C:269:LEU:HA	1.78	0.43
2:C:879:ARG:HD2	2:C:879:ARG:N	2.32	0.43
2:C:910:LYS:O	2:C:914:ILE:HG13	2.18	0.43
3:D:387:LEU:N	3:D:387:LEU:HD22	2.33	0.43
5:F:229:TYR:CE1	5:F:230:LYS:HD3	2.53	0.43
5:F:382:THR:OG1	5:F:385:GLU:HG3	2.18	0.43
2:C:10:ARG:NH2	2:C:12:VAL:HG12	2.33	0.43
2:C:405:ARG:HD2	2:C:442:GLU:OE2	2.19	0.43
2:C:774:LEU:HD22	5:F:350:LEU:CD1	2.49	0.43
3:D:248:PRO:HG3	3:D:308:LYS:HG3	1.99	0.43
3:D:473:LEU:HD21	3:D:495:ARG:NH2	2.32	0.43
3:D:1284:GLU:O	3:D:1284:GLU:HG3	2.17	0.43
4:E:54:LEU:O	4:E:55:PHE:HD2	2.01	0.43
2:C:137:VAL:HG12	2:C:138:SER:O	2.19	0.43
2:C:197:LEU:HD12	2:C:221:LEU:HD21	2.00	0.43
3:D:129:PHE:CZ	3:D:571:LYS:HB3	2.53	0.43
7:H:16:DC:H2''	7:H:17:DC:OP2	2.16	0.43
1:A:34:VAL:HG22	1:B:42:ARG:CZ	2.49	0.43
1:B:52:ALA:HB2	1:B:170:VAL:C	2.44	0.43
2:C:409:ARG:HD2	2:C:452:ILE:HG22	2.00	0.43
2:C:627:ARG:NH1	2:C:640:ARG:HG3	2.34	0.43
2:C:679:PHE:HA	3:D:943:THR:HG23	2.00	0.43
2:C:722:ILE:O	2:C:722:ILE:HG22	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:130:SER:HB2	5:F:83:GLN:HE22	1.82	0.43
3:D:1241:PHE:C	3:D:1243:THR:N	2.76	0.43
1:A:9:PRO:HD2	1:B:224:TYR:CD2	2.53	0.43
1:B:138:LEU:HD11	1:B:140:MET:HE3	2.01	0.43
2:C:460:ARG:HD2	2:C:485:TYR:CE1	2.54	0.43
3:D:534:ARG:HD2	5:F:312:GLN:OE1	2.18	0.43
3:D:1385:GLY:CA	3:D:1413:THR:HG23	2.48	0.43
3:D:1484:THR:O	4:E:25:LYS:NZ	2.36	0.43
5:F:274:THR:OG1	5:F:295:MET:HE1	2.18	0.43
1:A:101:LEU:HD11	1:A:109:VAL:CG1	2.49	0.43
2:C:278:GLU:HG2	2:C:284:ARG:HA	2.00	0.43
2:C:545:ASN:HB3	2:C:583:LEU:HD22	2.01	0.43
2:C:575:GLN:HG3	2:C:670:GLN:HG3	2.00	0.43
2:C:582:GLY:N	2:C:584:GLU:OE2	2.52	0.43
2:C:1057:SER:HB3	2:C:1058:ASP:H	1.38	0.43
3:D:122:GLU:HG2	3:D:152:LEU:HD11	2.00	0.43
3:D:371:ILE:HB	5:F:232:ARG:HH12	1.82	0.43
3:D:405:ASP:CB	3:D:423:ASP:OD1	2.66	0.43
3:D:576:GLU:OE2	5:F:80:PRO:HG3	2.19	0.43
1:B:65:PHE:CE2	3:D:809:PRO:HB2	2.53	0.43
2:C:486:MET:HE1	2:C:494:TYR:CD2	2.54	0.43
3:D:34:TYR:CE1	3:D:35:ARG:HG3	2.54	0.43
3:D:1152:GLU:HG2	3:D:1162:GLU:HB2	2.00	0.43
3:D:1197:ARG:HD2	3:D:1398:TRP:CE2	2.54	0.43
1:A:90:LEU:HD13	1:A:90:LEU:HA	1.66	0.43
1:B:99:LEU:HB2	1:B:142:VAL:HG22	2.00	0.43
2:C:144:PRO:CG	2:C:165:LEU:HD23	2.47	0.43
3:D:787:LEU:HD21	3:D:947:ILE:HG21	2.00	0.43
3:D:1066:THR:O	3:D:1067:VAL:C	2.62	0.43
3:D:1195:GLN:O	3:D:1398:TRP:HZ3	2.02	0.43
1:B:197:LEU:HD23	1:B:197:LEU:HA	1.64	0.43
2:C:226:VAL:O	2:C:229:MET:HG2	2.19	0.43
2:C:627:ARG:NE	2:C:639:GLN:O	2.52	0.43
3:D:879:ARG:HB3	3:D:902:LEU:CD1	2.48	0.43
3:D:893:GLU:H	3:D:894:LYS:HZ3	1.66	0.43
1:B:64:GLU:HA	1:B:165:ILE:HD13	2.01	0.42
1:B:101:LEU:HB2	1:B:114:PHE:CE1	2.54	0.42
2:C:202:TYR:OH	2:C:304:LEU:HB2	2.19	0.42
2:C:767:PRO:CB	2:C:771:GLU:HG2	2.49	0.42
2:C:939:ARG:HG2	2:C:939:ARG:H	1.40	0.42
2:C:1043:TYR:CG	3:D:763:MET:HG2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1066:ALA:HA	2:C:1077:PRO:HD2	2.01	0.42
3:D:536:ALA:HA	5:F:315:VAL:O	2.18	0.42
5:F:226:LYS:HB3	7:H:1:DT:H73	2.01	0.42
1:B:179:PHE:CB	1:B:197:LEU:HD23	2.48	0.42
2:C:258:TYR:CE2	2:C:293:PHE:HB2	2.54	0.42
2:C:436:GLY:HA2	2:C:538:GLN:O	2.19	0.42
2:C:673:LEU:HD12	2:C:673:LEU:HA	1.86	0.42
2:C:717:LEU:HD23	2:C:763:GLY:HA2	2.01	0.42
3:D:702:LEU:HD23	3:D:702:LEU:HA	1.65	0.42
3:D:810:GLU:H	3:D:810:GLU:CD	2.26	0.42
1:A:79:ILE:CD1	1:A:165:ILE:HG21	2.49	0.42
1:B:92:PRO:O	1:B:146:ARG:NH1	2.43	0.42
2:C:54:ILE:HG21	2:C:355:VAL:HG11	2.01	0.42
2:C:469:THR:HG23	2:C:471:TYR:CE1	2.54	0.42
3:D:313:MET:CE	3:D:318:ARG:C	2.92	0.42
2:C:708:TYR:HB3	2:C:790:LEU:HD21	2.00	0.42
3:D:462:GLN:HB2	3:D:513:ILE:HG21	2.01	0.42
3:D:844:ALA:O	3:D:867:ARG:HB3	2.20	0.42
3:D:879:ARG:HB3	3:D:902:LEU:HD11	2.01	0.42
3:D:1278:ASP:O	3:D:1294:VAL:HG12	2.19	0.42
2:C:247:PRO:HA	2:C:248:PRO:HD3	1.69	0.42
2:C:334:ARG:HB2	2:C:339:LEU:HD21	2.01	0.42
2:C:722:ILE:HD13	2:C:823:VAL:CG2	2.49	0.42
3:D:202:VAL:HG21	3:D:400:VAL:HG13	2.02	0.42
3:D:335:LEU:HD23	3:D:335:LEU:HA	1.81	0.42
3:D:411:THR:HB	3:D:437:VAL:H	1.85	0.42
3:D:711:LEU:HD13	3:D:778:LEU:HD13	2.01	0.42
1:A:57:TYR:CD1	1:A:161:ARG:HD2	2.55	0.42
1:B:131:THR:C	1:B:132:LEU:HD12	2.44	0.42
2:C:135:VAL:HG23	2:C:395:LYS:HG3	2.01	0.42
2:C:348:LEU:HA	2:C:348:LEU:HD12	1.79	0.42
2:C:391:LEU:HD23	2:C:391:LEU:HA	1.55	0.42
3:D:274:ARG:O	3:D:275:GLU:HB2	2.20	0.42
3:D:561:GLY:HA3	5:F:132:ARG:HD3	2.01	0.42
3:D:684:LYS:O	3:D:687:VAL:HG22	2.19	0.42
3:D:1296:SER:OG	3:D:1297:GLU:N	2.53	0.42
3:D:1493:LYS:HE3	3:D:1493:LYS:CA	2.34	0.42
1:A:206:THR:HG22	1:A:208:LEU:H	1.85	0.42
1:B:162:ILE:HG22	1:B:163:ASN:N	2.34	0.42
2:C:394:PHE:HE2	2:C:632:ASN:CB	2.19	0.42
2:C:875:GLY:O	2:C:879:ARG:HD3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1090:LYS:HA	2:C:1090:LYS:HD3	1.82	0.42
3:D:715:ALA:HB3	3:D:764:LEU:HA	2.00	0.42
3:D:1046:GLN:OE1	3:D:1247:ALA:HB2	2.19	0.42
3:D:1493:LYS:HD2	3:D:1493:LYS:N	2.34	0.42
2:C:694:LEU:HD23	2:C:694:LEU:HA	1.76	0.42
2:C:905:ILE:C	2:C:907:ASP:H	2.28	0.42
2:C:1104:GLU:H	2:C:1104:GLU:CD	2.28	0.42
1:A:228:PRO:HB3	1:B:13:VAL:HG21	2.01	0.42
1:B:86:VAL:HG12	1:B:123:MET:HG3	2.02	0.42
2:C:644:VAL:HG22	2:C:645:VAL:N	2.34	0.42
2:C:711:GLU:O	2:C:758:ARG:NH1	2.51	0.42
3:D:959:GLU:O	3:D:963:TYR:CD1	2.73	0.42
3:D:407:VAL:HA	3:D:422:ALA:CB	2.49	0.42
3:D:1045:MET:HE1	3:D:1245:GLY:C	2.45	0.42
3:D:1234:THR:CG2	3:D:1238:MET:HE3	2.48	0.42
5:F:315:VAL:HG22	5:F:316:SER:N	2.35	0.42
1:B:226:SER:O	1:B:228:PRO:HD3	2.19	0.41
2:C:174:LEU:HG	2:C:193:LEU:HD22	2.01	0.41
2:C:224:GLU:H	2:C:224:GLU:CD	2.28	0.41
2:C:229:MET:HB2	2:C:233:GLU:CB	2.40	0.41
2:C:424:GLY:O	2:C:428:ARG:HD2	2.19	0.41
3:D:210:ARG:HA	3:D:345:TYR:O	2.19	0.41
3:D:428:LYS:HE2	3:D:428:LYS:HB3	1.85	0.41
3:D:897:TRP:HH2	3:D:902:LEU:HD22	1.83	0.41
3:D:1314:LYS:O	3:D:1317:ASP:HB2	2.20	0.41
3:D:1493:LYS:HZ3	3:D:1496:GLU:HG3	1.83	0.41
4:E:46:PRO:HB2	4:E:57:ASP:HB3	2.00	0.41
6:G:12:DG:H5'	6:G:12:DG:C8	2.45	0.41
1:A:94:LEU:O	1:A:146:ARG:NH1	2.53	0.41
2:C:76:PRO:HA	2:C:77:PRO:HD3	1.93	0.41
2:C:129:ILE:HD11	2:C:386:PHE:CD2	2.55	0.41
2:C:926:PHE:HE1	2:C:929:ARG:NH1	2.17	0.41
3:D:23:TYR:CE2	3:D:89:ARG:HD3	2.54	0.41
3:D:474:GLU:HG3	3:D:496:LEU:HD11	2.02	0.41
3:D:957:PRO:HG3	3:D:1007:VAL:HA	2.02	0.41
3:D:1286:THR:HG22	3:D:1287:GLU:N	2.35	0.41
8:I:2:U:O2'	8:I:3:U:H5'	2.20	0.41
1:A:11:PHE:O	1:B:228:PRO:HA	2.20	0.41
2:C:127:PHE:O	2:C:133:ASP:HA	2.21	0.41
2:C:473:ARG:HA	2:C:531:PHE:CD2	2.55	0.41
2:C:906:PHE:CE1	3:D:1067:VAL:HA	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1001:VAL:CG1	3:D:630:VAL:HB	2.50	0.41
2:C:681:GLY:CA	3:D:939:PHE:CD2	3.03	0.41
3:D:501:ALA:CB	3:D:1452:ILE:HG22	2.48	0.41
1:B:80:LEU:HD21	3:D:842:VAL:HG12	2.02	0.41
3:D:90:MET:HE3	3:D:90:MET:HB3	1.87	0.41
3:D:172:PRO:HB2	3:D:175:VAL:CG2	2.50	0.41
3:D:313:MET:HE2	3:D:319:ALA:N	2.35	0.41
3:D:1099:VAL:HG13	3:D:1223:ILE:CD1	2.50	0.41
1:A:129:ILE:HD13	1:A:129:ILE:HA	1.84	0.41
1:B:206:THR:HG22	1:B:209:GLU:N	2.19	0.41
2:C:570:PRO:HD2	2:C:635:THR:HB	2.03	0.41
3:D:307:ALA:HB1	3:D:311:LEU:HD23	2.03	0.41
3:D:1489:GLN:OE1	3:D:1489:GLN:HA	2.21	0.41
5:F:148:LYS:HE3	5:F:148:LYS:HB2	1.84	0.41
8:I:0:A:C2'	8:I:1:A:O5'	2.69	0.41
1:A:228:PRO:HB3	1:B:13:VAL:CG2	2.50	0.41
2:C:328:LEU:HD23	2:C:328:LEU:HA	1.71	0.41
2:C:595:LEU:HD21	2:C:623:TYR:HB3	2.03	0.41
2:C:678:PRO:HA	2:C:683:ASN:HD22	1.86	0.41
2:C:684:PHE:O	2:C:685:GLU:C	2.64	0.41
3:D:625:TYR:CE1	3:D:751:LEU:HD11	2.55	0.41
3:D:832:ARG:HD3	3:D:832:ARG:HA	1.95	0.41
3:D:1028:ALA:O	3:D:1029:ARG:HG2	2.20	0.41
3:D:1312:LEU:HD21	3:D:1327:ARG:HG2	2.01	0.41
3:D:313:MET:HE2	3:D:318:ARG:C	2.46	0.41
3:D:500:ARG:HH12	3:D:1390:LEU:HD21	1.84	0.41
5:F:287:THR:HG22	5:F:288:TYR:N	2.34	0.41
7:H:18:DG:H1'	7:H:19:DG:H5'	2.02	0.41
1:A:225:PHE:O	1:B:11:PHE:CE2	2.74	0.41
2:C:12:VAL:HG21	2:C:472:ARG:HD3	2.02	0.41
2:C:55:GLU:O	2:C:56:GLU:CB	2.52	0.41
2:C:69:LEU:HD23	2:C:69:LEU:HA	1.90	0.41
2:C:161:SER:HA	2:C:172:ILE:O	2.21	0.41
2:C:235:LEU:HD21	2:C:254:VAL:CG2	2.51	0.41
2:C:501:THR:HA	2:C:502:PRO:HD3	1.96	0.41
2:C:540:PHE:HB3	2:C:544:THR:HB	2.03	0.41
2:C:775:ARG:HH21	5:F:422:LEU:CD2	2.34	0.41
3:D:84:ILE:HD12	3:D:87:ARG:HD2	2.01	0.41
3:D:701:LEU:O	3:D:702:LEU:HD23	2.20	0.41
3:D:893:GLU:N	3:D:894:LYS:HZ3	2.19	0.41
3:D:1197:ARG:HB2	3:D:1398:TRP:CH2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1240:THR:HG21	3:D:1251:ASP:OD2	2.20	0.41
1:B:123:MET:HB3	1:B:123:MET:HE3	1.85	0.41
1:B:124:ASN:HD22	1:B:127:LEU:HD22	1.83	0.41
2:C:586:ARG:HA	2:C:586:ARG:HD2	1.86	0.41
2:C:880:MET:HB3	3:D:1061:PHE:CD2	2.56	0.41
3:D:1069:GLU:O	3:D:1073:SER:OG	2.39	0.41
3:D:1205:TYR:CZ	3:D:1366:LYS:HD3	2.55	0.41
1:A:10:VAL:HG23	1:A:10:VAL:O	2.21	0.40
2:C:100:LEU:CD2	2:C:367:LEU:O	2.69	0.40
2:C:197:LEU:HD23	2:C:197:LEU:HA	1.93	0.40
2:C:577:PRO:HB3	2:C:993:PHE:CG	2.57	0.40
3:D:103:TRP:O	3:D:107:ASP:HB3	2.21	0.40
3:D:169:TYR:OH	3:D:198:ARG:HB2	2.21	0.40
3:D:186:VAL:HG12	3:D:187:LYS:N	2.35	0.40
3:D:689:ASP:OD2	4:E:51:LEU:HD11	2.21	0.40
3:D:1099:VAL:O	3:D:1103:HIS:HB3	2.21	0.40
5:F:408:LEU:HD23	5:F:408:LEU:HA	1.89	0.40
1:A:14:ARG:C	1:A:15:THR:CG2	2.94	0.40
1:A:115:LEU:HD23	1:A:115:LEU:HA	1.82	0.40
2:C:211:LEU:HD23	2:C:311:PHE:HD2	1.85	0.40
2:C:681:GLY:HA2	3:D:939:PHE:CE2	2.56	0.40
2:C:926:PHE:CE2	2:C:930:LYS:HD3	2.56	0.40
3:D:216:VAL:HB	3:D:382:GLU:HG2	2.02	0.40
3:D:237:LYS:O	3:D:240:GLU:HB2	2.20	0.40
3:D:693:GLU:CG	4:E:48:MET:HE1	2.51	0.40
5:F:364:ARG:NH1	5:F:364:ARG:HB3	2.37	0.40
1:A:18:ARG:O	1:A:207:PRO:HD3	2.21	0.40
2:C:64:LEU:N	2:C:103:LYS:HE2	2.36	0.40
2:C:778:PHE:CD1	5:F:422:LEU:HD11	2.57	0.40
3:D:465:LEU:HD23	3:D:465:LEU:HA	1.81	0.40
3:D:1141:GLU:CG	3:D:1168:MET:HE1	2.52	0.40
4:E:46:PRO:HB2	4:E:57:ASP:HB2	2.03	0.40
6:G:18:DT:C4	8:I:0:A:N6	2.77	0.40
1:B:80:LEU:HD23	1:B:80:LEU:HA	1.95	0.40
2:C:53:PRO:HB3	2:C:67:ASP:OD1	2.21	0.40
2:C:98:LEU:HA	2:C:98:LEU:HD23	1.81	0.40
2:C:773:LEU:O	2:C:777:ILE:HG13	2.21	0.40
3:D:185:VAL:HG13	3:D:189:GLN:HB3	2.02	0.40
3:D:372:ASP:OD1	3:D:372:ASP:N	2.54	0.40
3:D:405:ASP:CG	3:D:406:ASP:H	2.29	0.40
3:D:827:ILE:O	3:D:833:GLU:HA	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1143:GLY:O	3:D:1147:ARG:CD	2.58	0.40
2:C:1095:LEU:HD23	3:D:582:LEU:HD22	2.03	0.40
3:D:659:LYS:HA	3:D:659:LYS:HD2	1.95	0.40
3:D:1305:LEU:HD23	3:D:1305:LEU:N	2.36	0.40
3:D:1441:GLN:HE21	3:D:1441:GLN:HB3	1.56	0.40

All (2) 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)
3:D:985:ASP:N	3:D:1497:GLU:OE2[1_545]	2.09	0.11
3:D:34:TYR:OH	3:D:327:GLU:OE1[4_1359]	2.18	0.02

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	224/315 (71%)	221 (99%)	3 (1%)	0	100	100
1	B	220/315 (70%)	210 (96%)	9 (4%)	1 (0%)	24	53
2	C	1107/1119 (99%)	1078 (97%)	26 (2%)	3 (0%)	36	64
3	D	1498/1524 (98%)	1454 (97%)	39 (3%)	5 (0%)	36	64
4	E	92/99 (93%)	89 (97%)	2 (2%)	1 (1%)	11	35
5	F	344/423 (81%)	336 (98%)	5 (2%)	3 (1%)	14	40
All	All	3485/3795 (92%)	3388 (97%)	84 (2%)	13 (0%)	30	58

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	362	GLY
2	C	766	GLU

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Mol	Chain	Res	Type
3	D	1242	HIS
3	D	1249	ALA
5	F	322	GLY
5	F	323	ASP
5	F	145	PRO
1	B	154	GLU
3	D	1244	GLY
3	D	320	ALA
4	E	94	PRO
2	C	415	PRO
3	D	530	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	199/273 (73%)	187 (94%)	12 (6%)	17	45
1	B	195/273 (71%)	189 (97%)	6 (3%)	35	67
2	C	930/941 (99%)	877 (94%)	53 (6%)	18	48
3	D	1258/1279 (98%)	1172 (93%)	86 (7%)	14	40
4	E	83/88 (94%)	81 (98%)	2 (2%)	43	73
5	F	298/371 (80%)	288 (97%)	10 (3%)	32	65
All	All	2963/3225 (92%)	2794 (94%)	169 (6%)	18	48

All (169) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	34	VAL
1	A	97	VAL
1	A	104	GLU
1	A	112	ARG
1	A	122	ILE
1	A	133	GLU

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Mol	Chain	Res	Type
1	A	184	THR
1	A	186	LEU
1	A	189	ARG
1	A	206	THR
1	A	229	GLN
1	B	34	VAL
1	B	142	VAL
1	B	186	LEU
1	B	189	ARG
1	B	199	ILE
1	B	206	THR
2	C	8	ARG
2	C	15	LEU
2	C	56	GLU
2	C	81	ASP
2	C	103	LYS
2	C	107	LEU
2	C	177	GLU
2	C	182	VAL
2	C	194	VAL
2	C	196	LEU
2	C	218	VAL
2	C	221	LEU
2	C	251	ASP
2	C	261	ILE
2	C	265	ARG
2	C	331	ARG
2	C	336	VAL
2	C	342	ASP
2	C	358	ARG
2	C	364	GLU
2	C	377	PRO
2	C	427	VAL
2	C	429	ASP
2	C	449	ILE
2	C	454	SER
2	C	460	ARG
2	C	513	VAL
2	C	524	VAL
2	C	525	SER
2	C	575	GLN
2	C	578	VAL

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Mol	Chain	Res	Type
2	C	583	LEU
2	C	584	GLU
2	C	586	ARG
2	C	589	ARG
2	C	617	ASP
2	C	640	ARG
2	C	670	GLN
2	C	698	ASP
2	C	775	ARG
2	C	807	ARG
2	C	808	ARG
2	C	813	VAL
2	C	815	LEU
2	C	820	ARG
2	C	848	VAL
2	C	923	GLU
2	C	939	ARG
2	C	968	LEU
2	C	978	ARG
2	C	1001	VAL
2	C	1014	SER
2	C	1057	SER
3	D	30	GLU
3	D	40	GLU
3	D	81	THR
3	D	97	THR
3	D	106	LYS
3	D	114	THR
3	D	141	ILE
3	D	142	LEU
3	D	155	ASP
3	D	179	VAL
3	D	190	GLU
3	D	191	LEU
3	D	198	ARG
3	D	230	TRP
3	D	231	VAL
3	D	234	GLU
3	D	256	GLU
3	D	272	LEU
3	D	276	ASP
3	D	312	ARG

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Mol	Chain	Res	Type
3	D	325	GLU
3	D	355	VAL
3	D	360	ARG
3	D	362	GLU
3	D	372	ASP
3	D	387	LEU
3	D	411	THR
3	D	415	VAL
3	D	421	LEU
3	D	427	VAL
3	D	525	ARG
3	D	572	ARG
3	D	587	ARG
3	D	610	LYS
3	D	618	LEU
3	D	628	ARG
3	D	632	VAL
3	D	650	LEU
3	D	709	HIS
3	D	754	PHE
3	D	778	LEU
3	D	808	THR
3	D	810	GLU
3	D	817	GLU
3	D	832	ARG
3	D	864	VAL
3	D	875	THR
3	D	894	LYS
3	D	904	VAL
3	D	972	LEU
3	D	984	THR
3	D	995	LEU
3	D	1014	ASN
3	D	1039	CYS
3	D	1041	LEU
3	D	1044	LEU
3	D	1051	GLU
3	D	1062	ARG
3	D	1067	VAL
3	D	1073	SER
3	D	1155	VAL
3	D	1162	GLU

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Mol	Chain	Res	Type
3	D	1188	VAL
3	D	1195	GLN
3	D	1208	ASP
3	D	1219	GLU
3	D	1221	VAL
3	D	1252	ILE
3	D	1271	LYS
3	D	1283	ILE
3	D	1284	GLU
3	D	1288	GLU
3	D	1290	LEU
3	D	1301	LYS
3	D	1305	LEU
3	D	1307	LYS
3	D	1308	GLU
3	D	1313	VAL
3	D	1317	ASP
3	D	1351	GLU
3	D	1418	LYS
3	D	1470	ARG
3	D	1486	VAL
3	D	1493	LYS
3	D	1499	ARG
3	D	1500	LYS
4	E	50	THR
4	E	89	MET
5	F	120	THR
5	F	159	ILE
5	F	205	ARG
5	F	287	THR
5	F	338	LEU
5	F	347	GLN
5	F	363	GLU
5	F	364	ARG
5	F	377	ASP
5	F	415	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	38	ASN
1	B	63	HIS

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Mol	Chain	Res	Type
1	B	81	ASN
1	B	212	ASN
2	C	22	GLN
2	C	102	HIS
2	C	187	ASN
2	C	204	GLN
2	C	343	GLN
2	C	556	ASN
2	C	565	GLN
2	C	683	ASN
2	C	834	GLN
2	C	991	GLN
2	C	1047	HIS
3	D	294	HIS
3	D	316	GLN
3	D	442	ASN
3	D	463	GLN
3	D	560	GLN
3	D	611	GLN
3	D	640	HIS
3	D	724	GLN
3	D	855	HIS
3	D	991	GLN
3	D	1116	ASN
3	D	1323	GLN
3	D	1333	HIS
3	D	1359	GLN
4	E	49	GLN
5	F	83	GLN
5	F	90	GLN
5	F	170	HIS
5	F	218	GLN
5	F	402	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
8	I	3/4 (75%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 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
11	UTP	D	2005	9	29,30,30	2.97	11 (37%)	43,47,47	1.78	10 (23%)

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
11	UTP	D	2005	9	-	8/22/38/38	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	D	2005	UTP	O4'-C1'	8.87	1.62	1.42
11	D	2005	UTP	O4'-C4'	-7.12	1.29	1.45
11	D	2005	UTP	C2'-C1'	-6.69	1.32	1.53
11	D	2005	UTP	O2'-C2'	3.77	1.52	1.43
11	D	2005	UTP	PB-O3A	3.70	1.63	1.59
11	D	2005	UTP	C5'-C4'	3.08	1.60	1.51
11	D	2005	UTP	C5-C4	-2.93	1.37	1.43
11	D	2005	UTP	PA-O5'	2.33	1.68	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	D	2005	UTP	O4-C4	-2.15	1.20	1.24
11	D	2005	UTP	C2-N1	-2.08	1.35	1.38
11	D	2005	UTP	C4-N3	-2.02	1.35	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	D	2005	UTP	N3-C2-N1	5.22	121.68	114.89
11	D	2005	UTP	C4-N3-C2	-4.91	120.51	126.61
11	D	2005	UTP	O4-C4-C5	-3.30	119.47	125.16
11	D	2005	UTP	C5-C4-N3	3.21	119.30	114.80
11	D	2005	UTP	O2'-C2'-C3'	2.61	120.17	111.82
11	D	2005	UTP	C2'-C1'-N1	-2.34	106.74	113.25
11	D	2005	UTP	C1'-N1-C2	2.34	121.79	117.59
11	D	2005	UTP	C2'-C3'-C4'	2.30	107.05	102.61
11	D	2005	UTP	C3'-C2'-C1'	2.23	105.68	101.46
11	D	2005	UTP	O2-C2-N1	-2.06	120.12	122.80

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	D	2005	UTP	C5'-O5'-PA-O1A
11	D	2005	UTP	C5'-O5'-PA-O3A
11	D	2005	UTP	C3'-C4'-C5'-O5'
11	D	2005	UTP	O4'-C4'-C5'-O5'
11	D	2005	UTP	C5'-O5'-PA-O2A
11	D	2005	UTP	C4'-C5'-O5'-PA
11	D	2005	UTP	PB-O3A-PA-O2A
11	D	2005	UTP	PB-O3A-PA-O1A

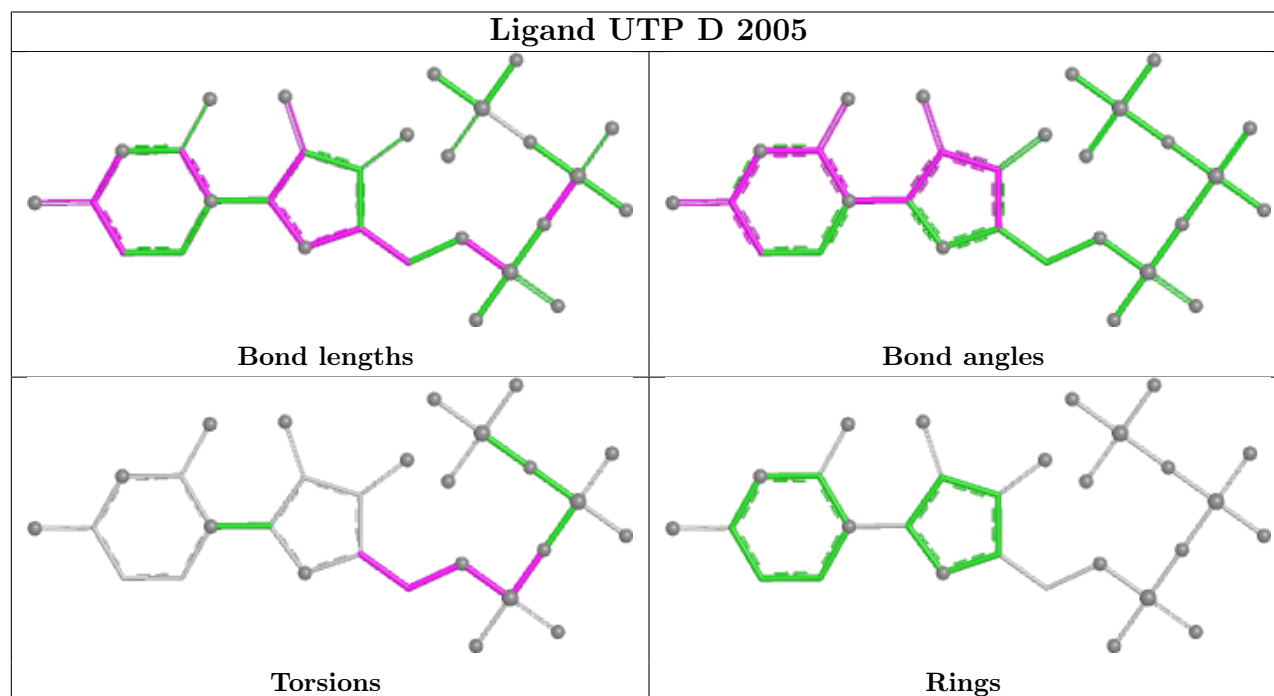
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	D	2005	UTP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	226/315 (71%)	-0.19	0 100 100	54, 77, 100, 109	0
1	B	222/315 (70%)	-0.23	1 (0%) 87 83	52, 77, 106, 122	0
2	C	1111/1119 (99%)	-0.04	13 (1%) 76 69	38, 80, 149, 164	0
3	D	1500/1524 (98%)	-0.21	13 (0%) 81 74	36, 73, 127, 171	0
4	E	94/99 (94%)	0.15	3 (3%) 50 40	48, 85, 133, 147	0
5	F	346/423 (81%)	-0.01	6 (1%) 69 60	49, 93, 137, 149	0
6	G	15/21 (71%)	-0.33	1 (6%) 24 19	49, 77, 152, 153	0
7	H	16/27 (59%)	-0.33	0 100 100	73, 99, 129, 129	0
8	I	4/4 (100%)	-0.30	0 100 100	56, 62, 80, 107	0
All	All	3534/3847 (91%)	-0.13	37 (1%) 79 72	36, 78, 138, 171	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
6	G	20	DT	4.2
3	D	1241	PHE	3.5
2	C	207	LEU	3.2
5	F	141	VAL	3.2
2	C	107	LEU	3.1
2	C	242	LEU	2.9
3	D	1244	GLY	2.8
3	D	1245	GLY	2.8
3	D	1242	HIS	2.8
2	C	196	LEU	2.8
1	B	228	PRO	2.7
3	D	412	GLY	2.6
3	D	1494	ALA	2.6
5	F	349	LEU	2.5
3	D	1128	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
5	F	323	ASP	2.4
2	C	394	PHE	2.3
2	C	421	GLU	2.3
2	C	101	ILE	2.3
2	C	171	TRP	2.3
4	E	32	ARG	2.3
3	D	420	VAL	2.3
3	D	1250	ALA	2.2
2	C	264	PRO	2.2
5	F	322	GLY	2.2
4	E	2	ALA	2.2
3	D	1486	VAL	2.2
3	D	1243	THR	2.2
2	C	102	HIS	2.2
2	C	186	VAL	2.2
3	D	1313	VAL	2.2
2	C	66	LEU	2.2
5	F	324	GLU	2.2
5	F	142	ARG	2.1
2	C	51	THR	2.1
4	E	93	TYR	2.1
3	D	1236	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

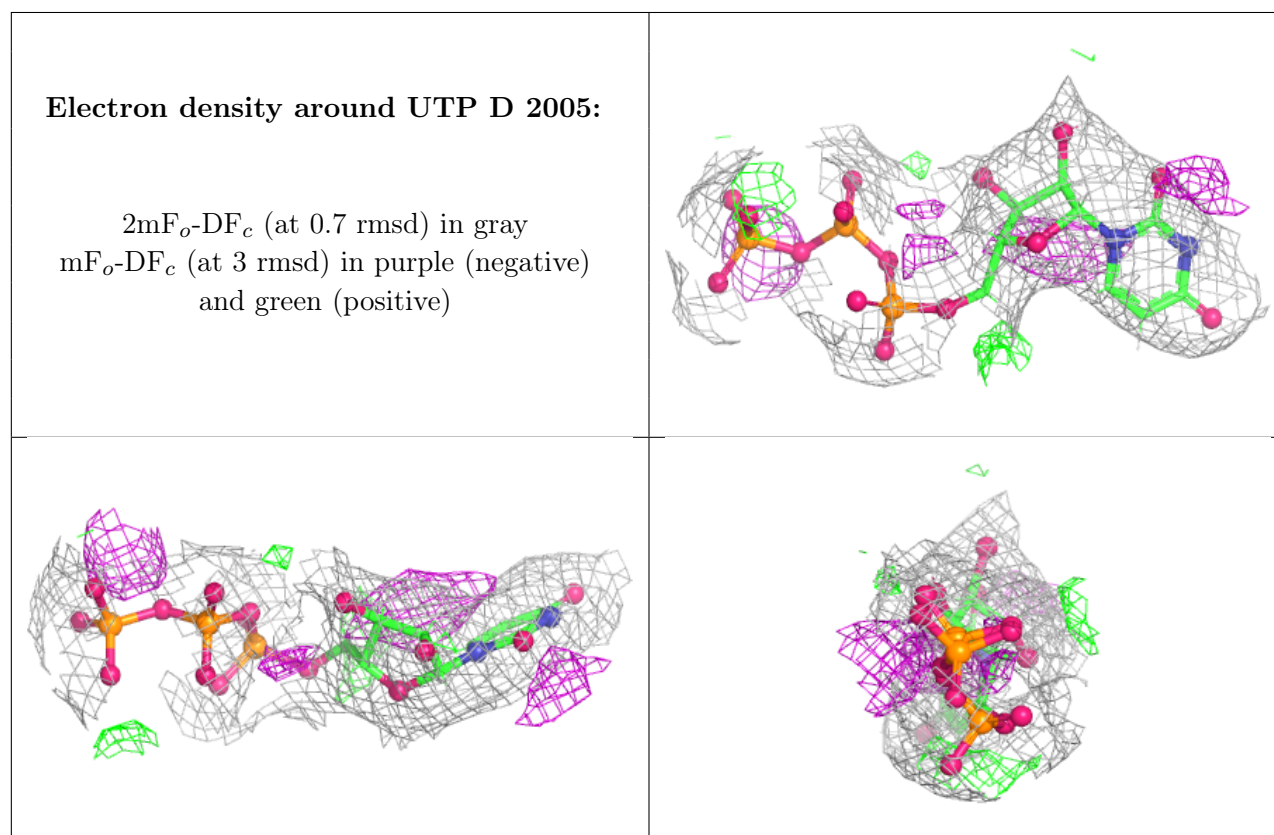
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
9	MG	D	2002	1/1	0.81	0.15	65,65,65,65	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
11	UTP	D	2005	29/29	0.86	0.09	53,71,98,110	0
9	MG	D	2001	1/1	0.97	0.04	40,40,40,40	0
10	ZN	D	2004	1/1	0.99	0.03	103,103,103,103	0
10	ZN	D	2003	1/1	1.00	0.07	66,66,66,66	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers ⓘ

There are no such residues in this entry.