



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:43 PM GMT

PDB ID : 1UU1
Title : COMPLEX OF HISTIDINOL-PHOSPHATE AMINOTRANSFERASE
(HISC) FROM THERMOTOGA MARITIMA (APO-FORM)
Authors : Vega, M.C.; Fernandez, F.J.; Lehman, F.; Wilmanns, M.
Deposited on : 2003-12-12
Resolution : 2.38 Å(reported)

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

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

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

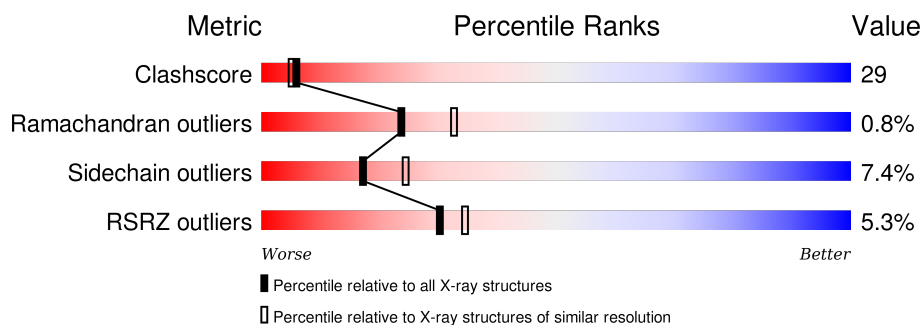
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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))
Clashscore	102246	4595 (2.40-2.36)
Ramachandran outliers	100387	4520 (2.40-2.36)
Sidechain outliers	100360	4522 (2.40-2.36)
RSRZ outliers	91569	4034 (2.40-2.36)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	335	<div> <div>6%</div> <div>56%39%</div> <div>..</div> </div>
1	B	335	<div> <div>3%</div> <div>56%40%</div> <div>.</div> </div>
1	C	335	<div> <div>6%</div> <div>54%41%5%</div> <div></div> </div>
1	D	335	<div> <div>7%</div> <div>50%45%</div> <div>..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	HSA	A	1336	X	-	-	-
3	HSA	B	1336	X	-	-	X
3	HSA	C	1336	X	-	-	-
3	HSA	D	1336	X	-	-	-

2 Entry composition [i](#)

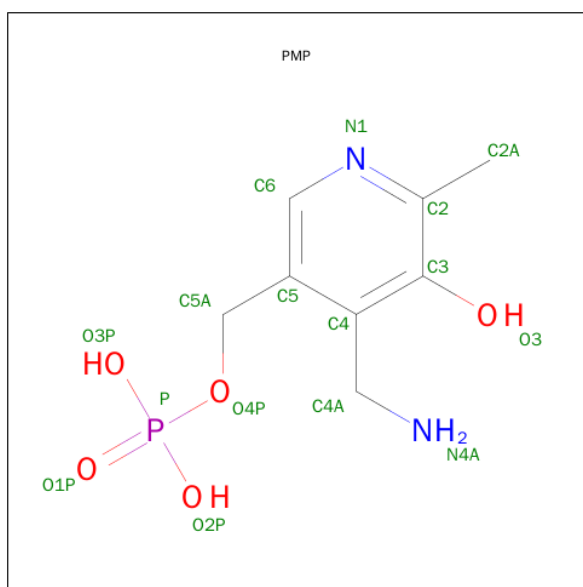
There are 4 unique types of molecules in this entry. The entry contains 11516 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HISTIDINOL-PHOSPHATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	1
			2727	1754	460	504	9			
1	B	335	Total	C	N	O	S	1	0	1
			2766	1778	466	512	10			
1	C	335	Total	C	N	O	S	0	0	0
			2772	1781	467	514	10			
1	D	335	Total	C	N	O	S	3	0	1
			2766	1778	466	512	10			

- Molecule 2 is 4'-DEOXY-4'-AMINOPYRIDOXAL-5'-PHOSPHATE (three-letter code: PMP) (formula: C₈H₁₃N₂O₅P).



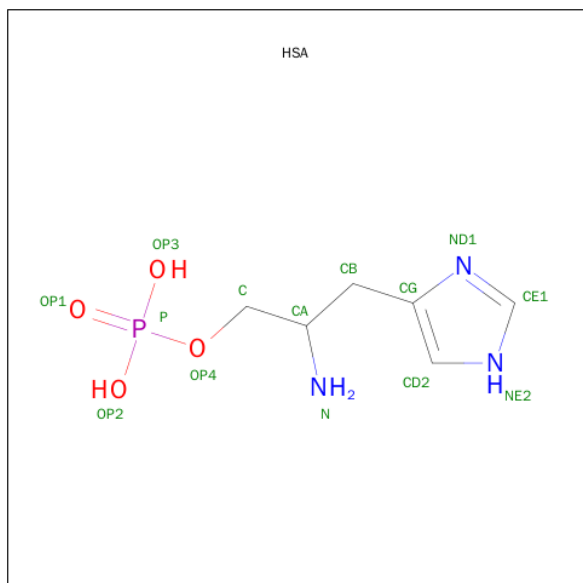
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
2	B	1	Total	C	N	O	P	0	0
			16	8	2	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
2	D	1	Total	C	N	O	P	0	0
			16	8	2	5	1		

- Molecule 3 is PHOSPHORIC ACID MONO-[2-AMINO-3-(3H-IMIDAZOL-4-YL)-PROPYL]ESTER (three-letter code: HSA) (formula: $C_6H_{12}N_3O_4P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	1	0
			13	6	2	4	1		
3	B	1	Total	C	N	O	P	0	0
			13	6	2	4	1		
3	C	1	Total	C	N	O	P	0	0
			13	6	2	4	1		
3	D	1	Total	C	N	O	P	2	0
			13	6	2	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	102	Total	O	0	0
			102	102		
4	B	82	Total	O	0	0
			82	82		
4	C	97	Total	O	0	0
			97	97		

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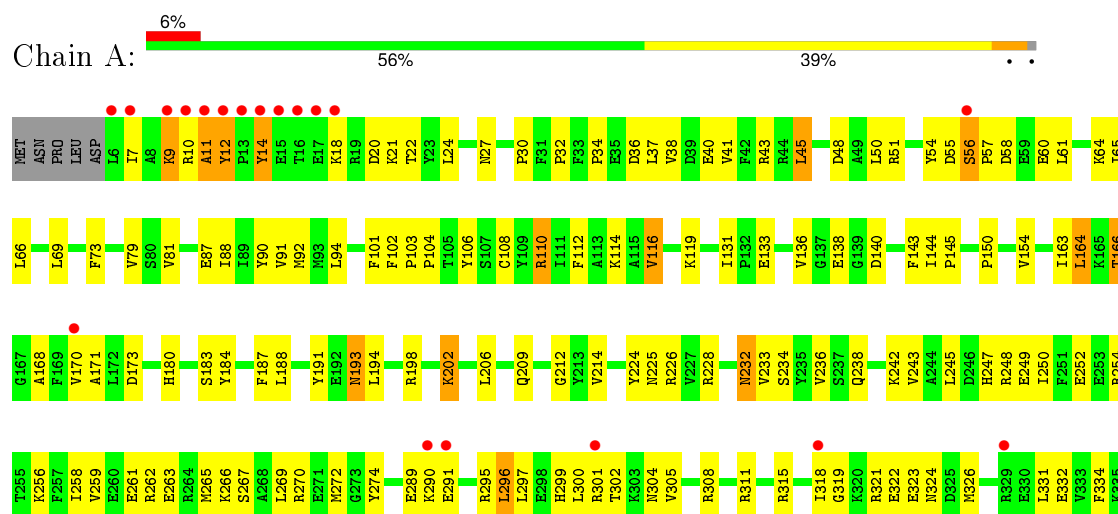
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	88	Total	O	0	0
			88	88		

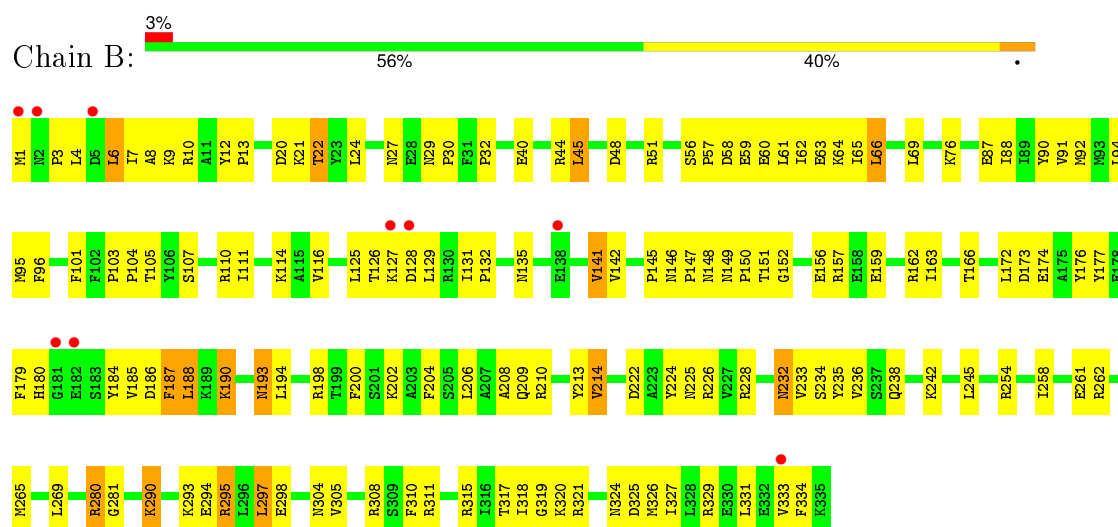
3 Residue-property plots

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

• Molecule 1: HISTIDINOL-PHOSPHATE AMINOTRANSFERASE

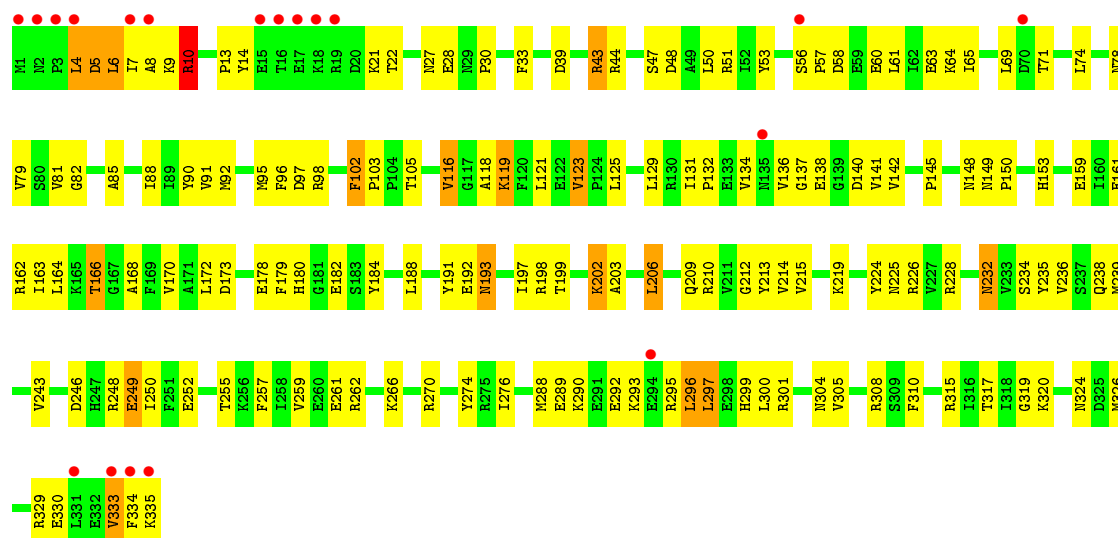


• Molecule 1: HISTIDINOL-PHOSPHATE AMINOTRANSFERASE

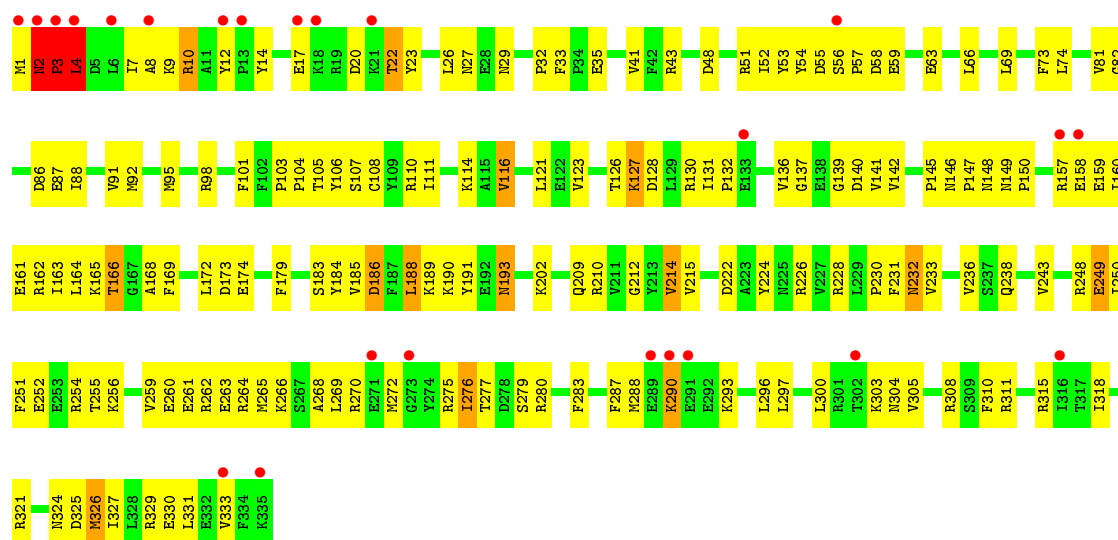


• Molecule 1: HISTIDINOL-PHOSPHATE AMINOTRANSFERASE





● Molecule 1: HISTIDINOL-PHOSPHATE AMINOTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.61Å 136.52Å 92.81Å 90.00° 95.19° 90.00°	Depositor
Resolution (Å)	19.67 – 2.38 19.67 – 2.35	Depositor EDS
% Data completeness (in resolution range)	90.7 (19.67-2.38) 89.8 (19.67-2.35)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.40 (at 2.35Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.206 , 0.268 0.210 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.411	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 49533 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11516	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HSA, PMP

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	0.57	0/2786	0.72	1/3757 (0.0%)
1	B	0.57	0/2826	0.72	0/3812
1	C	0.53	0/2832	0.71	1/3817 (0.0%)
1	D	0.53	0/2826	0.69	1/3812 (0.0%)
All	All	0.55	0/11270	0.71	3/15198 (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
1	A	0	1
1	B	0	2
1	C	0	1
All	All	0	4

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	5	ASP	N-CA-C	-5.95	94.93	111.00
1	D	4	LEU	N-CA-C	-5.47	96.23	111.00
1	A	56	SER	C-N-CA	-5.46	99.08	122.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	90	TYR	Sidechain
1	B	176	TYR	Sidechain
1	B	90	TYR	Sidechain
1	C	90	TYR	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	2727	0	2702	146	0
1	B	2766	0	2742	164	0
1	C	2772	0	2746	188	0
1	D	2766	0	2742	185	0
2	A	16	0	9	4	0
2	B	16	0	8	4	0
2	C	16	0	9	3	0
2	D	16	0	8	5	0
3	A	13	0	7	1	0
3	B	13	0	7	0	0
3	C	13	0	7	3	0
3	D	13	0	7	1	0
4	A	102	0	0	9	0
4	B	82	0	0	7	0
4	C	97	0	0	8	0
4	D	88	0	0	9	0
All	All	11516	0	10994	632	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 29.

All (632) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:THR:HG22	1:D:128:ASP:H	1.07	1.11
1:C:43:ARG:HG2	1:C:43:ARG:HH21	1.24	1.02
1:C:58:ASP:HB2	1:C:238:GLN:HE22	1.23	1.01
1:C:91:VAL:HG21	1:C:224:TYR:CE1	1.97	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:LYS:NZ	2:A:1335:PMP:H4A2	1.78	0.99
1:C:71:THR:HG21	1:C:74:LEU:HB2	1.45	0.98
1:C:193:ASN:HD22	1:C:193:ASN:H	0.97	0.95
1:B:24:LEU:HD13	1:B:327:ILE:HD11	1.45	0.95
1:D:58:ASP:HB2	1:D:238:GLN:NE2	1.80	0.93
1:C:58:ASP:HB2	1:C:238:GLN:NE2	1.83	0.92
1:C:22:THR:HG23	1:C:305:VAL:HA	1.51	0.92
1:C:232:ASN:H	1:C:232:ASN:HD22	1.10	0.92
1:B:58:ASP:HB2	1:B:238:GLN:NE2	1.85	0.90
1:A:193:ASN:H	1:A:193:ASN:HD22	1.17	0.90
1:D:232:ASN:HD22	1:D:232:ASN:H	1.19	0.90
1:D:88:ILE:HD13	1:D:214:VAL:HG22	1.54	0.88
1:B:88:ILE:HD13	1:B:214:VAL:HG22	1.52	0.88
1:B:92:MET:HA	1:B:95:MET:CE	2.05	0.87
1:D:56:SER:HB3	1:D:57:PRO:HD3	1.57	0.87
1:D:126:THR:HG22	1:D:128:ASP:N	1.90	0.87
1:D:290:LYS:H	1:D:290:LYS:HD2	1.40	0.86
1:C:91:VAL:HG21	1:C:224:TYR:HE1	1.40	0.86
1:C:119:LYS:CD	1:C:119:LYS:H	1.89	0.86
1:D:188:LEU:HG	4:D:2052:HOH:O	1.73	0.85
1:C:193:ASN:ND2	1:C:193:ASN:H	1.75	0.85
1:A:91:VAL:HG21	1:A:224:TYR:HE1	1.40	0.84
1:C:261:GLU:HG2	1:C:324:ASN:ND2	1.92	0.84
1:A:166:THR:HG22	1:A:168:ALA:H	1.41	0.84
1:B:92:MET:HA	1:B:95:MET:HE3	1.60	0.83
1:D:193:ASN:HD22	1:D:193:ASN:H	1.25	0.83
1:A:236:VAL:H	1:B:209:GLN:HE22	1.25	0.83
1:B:232:ASN:H	1:B:232:ASN:HD22	1.25	0.83
1:C:56:SER:HB3	1:C:57:PRO:HD3	1.59	0.83
1:A:88:ILE:HD13	1:A:214:VAL:HG22	1.57	0.83
1:C:288:MET:HE3	1:C:293:LYS:HA	1.62	0.81
1:D:58:ASP:HB2	1:D:238:GLN:HE22	1.42	0.80
1:C:193:ASN:HD22	1:C:193:ASN:N	1.79	0.80
1:B:24:LEU:CD1	1:B:327:ILE:HD11	2.11	0.79
1:A:272:MET:HE3	1:A:332:GLU:HA	1.64	0.79
1:B:193:ASN:H	1:B:193:ASN:HD22	1.30	0.79
1:A:91:VAL:HG21	1:A:224:TYR:CE1	2.17	0.79
1:A:48:ASP:HA	1:B:32:PRO:HB3	1.65	0.78
1:A:30:PRO:HB3	1:A:319:GLY:HA2	1.66	0.77
1:C:232:ASN:N	1:C:232:ASN:HD22	1.78	0.77
1:A:94:LEU:HD21	1:A:116:VAL:HG22	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:LYS:HZ2	2:A:1335:PMP:H4A2	1.49	0.76
1:B:94:LEU:CD2	1:B:116:VAL:HG11	2.16	0.76
1:C:232:ASN:ND2	1:C:232:ASN:H	1.83	0.75
1:B:107:SER:O	1:B:111:ILE:HD13	1.86	0.74
1:C:225:ASN:HD22	1:C:228:ARG:HE	1.33	0.74
1:D:248:ARG:HA	1:D:251:PHE:HD1	1.52	0.74
1:C:119:LYS:HD3	1:C:119:LYS:H	1.52	0.73
1:C:225:ASN:ND2	1:C:228:ARG:HE	1.86	0.73
1:B:96:PHE:CG	1:B:141:VAL:HG22	2.23	0.73
1:C:226:ARG:NE	1:D:8:ALA:HA	2.03	0.73
1:D:270:ARG:HE	1:D:276:ILE:HD11	1.53	0.73
1:A:131:ILE:HG21	1:A:163:ILE:HD11	1.70	0.73
1:C:209:GLN:HE22	1:D:236:VAL:H	1.37	0.72
1:B:225:ASN:ND2	1:B:228:ARG:HE	1.87	0.72
1:D:215:VAL:HG12	4:D:2052:HOH:O	1.89	0.72
1:D:7:ILE:C	1:D:9:LYS:H	1.93	0.72
1:D:91:VAL:HG21	1:D:224:TYR:CE1	2.25	0.71
1:B:22:THR:HG23	1:B:305:VAL:HA	1.71	0.71
1:A:9:LYS:HG3	1:B:226:ARG:CZ	2.21	0.71
1:B:311:ARG:HH22	1:D:249:GLU:CG	2.03	0.71
1:D:290:LYS:N	1:D:290:LYS:HD2	2.05	0.71
1:C:145:PRO:HA	1:C:173:ASP:HB3	1.72	0.70
1:B:131:ILE:HG21	1:B:163:ILE:HD11	1.71	0.70
1:A:295:ARG:HB3	1:A:334:PHE:HZ	1.56	0.70
1:C:142:VAL:HG11	1:C:163:ILE:HG21	1.72	0.70
1:D:2:ASN:O	1:D:4:LEU:N	2.25	0.70
1:D:131:ILE:HG21	1:D:163:ILE:HD11	1.74	0.70
1:C:102:PHE:HB3	1:C:123:VAL:O	1.92	0.70
1:C:58:ASP:CB	1:C:238:GLN:HE22	2.02	0.70
1:A:65:ILE:O	1:A:69:LEU:HD13	1.91	0.69
1:A:21:LYS:HG2	1:A:304:ASN:HB3	1.74	0.69
1:A:9:LYS:HB2	1:B:226:ARG:HD2	1.74	0.69
1:B:311:ARG:HH12	1:D:249:GLU:HG2	1.57	0.69
1:A:225:ASN:ND2	1:A:228:ARG:HE	1.89	0.69
1:D:4:LEU:CD1	1:D:7:ILE:HD12	2.23	0.69
1:A:247:HIS:HB3	1:A:250:ILE:HD13	1.75	0.69
1:C:202:LYS:HE3	2:C:1335:PMP:H4A2	1.73	0.69
1:C:43:ARG:HG2	1:C:43:ARG:NH2	2.03	0.69
1:C:6:LEU:HD13	1:C:7:ILE:N	2.08	0.69
1:A:249:GLU:HB2	4:A:2085:HOH:O	1.91	0.68
1:A:225:ASN:HD22	1:A:228:ARG:HE	1.39	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:SER:HB3	1:B:57:PRO:HD3	1.74	0.68
1:B:96:PHE:CB	1:B:141:VAL:HG22	2.23	0.68
1:A:299:HIS:O	1:A:302:THR:HB	1.93	0.68
1:D:57:PRO:HG2	1:D:228:ARG:NH2	2.07	0.68
1:B:22:THR:HG22	4:B:2070:HOH:O	1.91	0.68
1:A:145:PRO:HA	1:A:173:ASP:HB3	1.76	0.68
1:A:57:PRO:HG3	1:A:81:VAL:HG12	1.74	0.68
1:A:209:GLN:HE22	1:B:236:VAL:H	1.42	0.67
1:A:266:LYS:O	1:A:270:ARG:HG3	1.93	0.67
1:C:119:LYS:H	1:C:119:LYS:CE	2.08	0.67
1:C:164:LEU:HG	1:C:191:TYR:CD2	2.29	0.67
1:B:232:ASN:HD22	1:B:232:ASN:N	1.91	0.67
1:C:95:MET:O	1:D:4:LEU:HD21	1.94	0.67
1:A:193:ASN:H	1:A:193:ASN:ND2	1.91	0.67
1:A:166:THR:CG2	1:A:168:ALA:H	2.08	0.67
1:B:94:LEU:HD23	1:B:116:VAL:HG11	1.76	0.67
1:D:142:VAL:HG11	1:D:163:ILE:HG21	1.76	0.66
1:C:163:ILE:O	1:C:166:THR:HB	1.95	0.66
1:A:136:VAL:HG23	1:A:166:THR:HG21	1.77	0.66
1:A:58:ASP:HB2	1:A:238:GLN:HE22	1.60	0.66
1:A:202:LYS:HZ1	2:A:1335:PMP:H4A2	1.57	0.66
1:C:166:THR:HG22	1:C:168:ALA:H	1.59	0.66
1:C:88:ILE:HG21	1:C:197:ILE:HD12	1.78	0.66
1:B:92:MET:HA	1:B:95:MET:HE2	1.77	0.66
1:B:224:TYR:CE2	1:B:228:ARG:HD3	2.30	0.65
1:B:96:PHE:HB2	1:B:141:VAL:HG22	1.79	0.65
1:B:22:THR:HG21	1:B:305:VAL:HG22	1.77	0.65
1:A:272:MET:CE	1:A:332:GLU:HA	2.26	0.65
1:B:329:ARG:O	1:B:333:VAL:HG22	1.97	0.65
1:C:92:MET:HE3	1:C:197:ILE:HD11	1.79	0.64
1:D:17:GLU:HG3	1:D:26:LEU:HD11	1.79	0.64
1:C:159:GLU:HA	1:C:162:ARG:NH2	2.12	0.64
1:B:58:ASP:HB2	1:B:238:GLN:HE22	1.63	0.64
1:B:327:ILE:O	1:B:331:LEU:HG	1.98	0.64
1:C:119:LYS:H	1:C:119:LYS:HE2	1.63	0.64
1:A:232:ASN:N	1:A:232:ASN:HD22	1.96	0.64
1:C:219:LYS:HB3	1:C:219:LYS:NZ	2.13	0.64
1:C:91:VAL:HG21	1:C:224:TYR:CD1	2.33	0.63
1:D:276:ILE:O	1:D:276:ILE:HD13	1.98	0.63
1:B:315:ARG:HD2	4:B:2076:HOH:O	1.97	0.63
1:C:58:ASP:CB	1:C:238:GLN:NE2	2.58	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:236:VAL:H	1:D:209:GLN:HE22	1.47	0.63
1:A:61:LEU:O	1:A:65:ILE:HG12	1.97	0.63
1:A:27:ASN:HD21	1:A:315:ARG:HD3	1.64	0.63
1:D:232:ASN:HD22	1:D:232:ASN:N	1.92	0.63
1:C:91:VAL:CG2	1:C:224:TYR:HE1	2.11	0.62
1:D:7:ILE:O	1:D:8:ALA:HB3	1.99	0.62
1:D:315:ARG:HD2	4:D:2084:HOH:O	1.99	0.62
1:D:264:ARG:NH2	1:D:325:ASP:OD2	2.33	0.62
1:B:63:GLU:HG3	1:B:76:LYS:HE2	1.81	0.62
1:B:64:LYS:HD3	1:B:242:LYS:HG2	1.82	0.62
1:C:136:VAL:HG12	1:C:137:GLY:N	2.14	0.62
1:A:296:LEU:HD22	1:A:300:LEU:HD11	1.80	0.62
1:D:51:ARG:HG3	1:D:52:ILE:HG23	1.80	0.62
1:A:248:ARG:O	1:A:252:GLU:HG3	2.00	0.62
1:A:56:SER:HB3	1:A:57:PRO:HD3	1.81	0.62
1:B:148:ASN:HB3	1:B:151:THR:OG1	2.00	0.62
1:D:290:LYS:H	1:D:290:LYS:CD	2.03	0.62
1:B:94:LEU:HD21	1:B:116:VAL:HG11	1.82	0.62
1:D:27:ASN:HD21	1:D:315:ARG:HD3	1.65	0.62
1:C:98:ARG:HA	1:C:119:LYS:HG2	1.81	0.62
1:B:24:LEU:HD13	1:B:327:ILE:CD1	2.26	0.61
1:C:98:ARG:HB3	1:C:140:ASP:OD2	1.99	0.61
1:C:88:ILE:HD12	1:C:214:VAL:CG2	2.30	0.61
1:C:136:VAL:HG12	1:C:137:GLY:H	1.65	0.61
1:C:137:GLY:O	1:C:140:ASP:HB2	2.01	0.61
1:D:185:VAL:O	1:D:188:LEU:HB2	2.00	0.61
1:B:311:ARG:HH22	1:D:249:GLU:HG2	1.65	0.61
1:C:57:PRO:HD2	1:D:14:TYR:OH	2.01	0.61
1:B:126:THR:HG22	1:B:127:LYS:N	2.16	0.60
1:D:86:ASP:OD1	2:D:1335:PMP:H5A2	2.01	0.60
1:A:323:GLU:O	1:A:326:MET:HG3	2.00	0.60
1:A:10:ARG:HG3	4:A:2003:HOH:O	1.99	0.60
1:B:333:VAL:O	1:B:333:VAL:HG23	2.01	0.60
1:D:29:ASN:HB2	1:D:202:LYS:O	2.00	0.60
1:C:105:THR:HA	1:C:149:ASN:O	2.01	0.60
1:C:192:GLU:HG3	1:D:1:MET:HG3	1.82	0.60
1:A:322:GLU:CD	1:A:322:GLU:H	2.02	0.60
1:B:311:ARG:NH1	1:D:249:GLU:HG2	2.17	0.60
1:C:85:ALA:HB3	2:C:1335:PMP:H5A2	1.82	0.60
1:C:88:ILE:HD12	1:C:214:VAL:HG22	1.84	0.60
1:B:60:GLU:HG3	1:B:242:LYS:HD3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:PRO:HG2	1:D:148:ASN:OD1	2.02	0.59
1:D:330:GLU:O	1:D:333:VAL:HG12	2.02	0.59
1:C:234:SER:O	1:C:238:GLN:HG3	2.02	0.59
1:C:57:PRO:HG3	1:C:81:VAL:HG12	1.84	0.59
1:A:58:ASP:CB	1:A:238:GLN:HE22	2.15	0.59
1:B:172:LEU:CD2	1:B:184:TYR:HB2	2.32	0.59
1:C:131:ILE:HG21	1:C:163:ILE:HD11	1.83	0.59
1:D:248:ARG:HA	1:D:251:PHE:CD1	2.37	0.59
1:C:180:HIS:CE1	1:C:182:GLU:HB3	2.38	0.59
1:B:145:PRO:HA	1:B:173:ASP:HB3	1.83	0.59
1:A:87:GLU:O	1:A:91:VAL:HG22	2.02	0.58
1:C:210:ARG:HD3	1:D:231:PHE:O	2.03	0.58
1:A:261:GLU:HG2	1:A:324:ASN:ND2	2.18	0.58
1:C:56:SER:O	1:C:58:ASP:N	2.36	0.58
1:B:172:LEU:HD22	1:B:184:TYR:HB2	1.85	0.58
1:A:188:LEU:HD22	1:A:194:LEU:HD23	1.85	0.58
1:C:121:LEU:HD12	1:C:123:VAL:HG22	1.84	0.58
1:A:315:ARG:HH22	3:A:1336:HSA:P	2.26	0.58
1:C:329:ARG:O	1:C:333:VAL:HG22	2.04	0.58
1:C:97:ASP:O	1:C:119:LYS:HG3	2.04	0.58
1:A:248:ARG:HD3	4:A:2086:HOH:O	2.03	0.58
1:A:58:ASP:H	1:A:238:GLN:HE22	1.51	0.58
1:B:116:VAL:O	1:B:116:VAL:HG12	2.03	0.58
1:C:43:ARG:HH21	1:C:43:ARG:CG	2.06	0.57
1:B:126:THR:HG22	1:B:127:LYS:H	1.69	0.57
1:A:133:GLU:CD	1:A:133:GLU:H	2.06	0.57
1:D:293:LYS:HD3	4:D:2083:HOH:O	2.04	0.57
1:D:4:LEU:HD11	1:D:7:ILE:HD12	1.86	0.57
1:C:57:PRO:HG2	1:C:228:ARG:NH2	2.19	0.57
1:D:303:LYS:O	1:D:305:VAL:HG23	2.04	0.57
1:D:174:GLU:OE2	1:D:183:SER:HA	2.05	0.57
1:A:11:ALA:HA	1:B:226:ARG:NH2	2.20	0.57
1:D:107:SER:O	1:D:111:ILE:HG12	2.04	0.57
1:A:88:ILE:O	1:A:92:MET:HG3	2.04	0.57
1:A:226:ARG:HG2	1:B:12:TYR:O	2.05	0.57
1:D:261:GLU:HG2	1:D:324:ASN:ND2	2.19	0.57
1:C:65:ILE:O	1:C:69:LEU:HD23	2.04	0.57
1:D:232:ASN:ND2	1:D:232:ASN:H	1.97	0.57
1:A:236:VAL:HG23	1:B:209:GLN:NE2	2.19	0.57
1:D:202:LYS:NZ	2:D:1335:PMP:H4A2	2.20	0.56
1:C:22:THR:HG21	1:C:305:VAL:HG22	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:ALA:HA	1:B:226:ARG:HH22	1.69	0.56
1:B:96:PHE:HB2	1:B:141:VAL:CG2	2.34	0.56
1:D:308:ARG:HG3	1:D:310:PHE:CE2	2.40	0.56
1:B:295:ARG:HH21	1:B:295:ARG:HG2	1.69	0.56
1:A:272:MET:HE2	1:A:274:TYR:HE1	1.70	0.56
1:B:185:VAL:O	1:B:188:LEU:HB2	2.05	0.56
1:C:178:GLU:HG2	1:C:203:ALA:CB	2.35	0.56
1:B:254:ARG:O	1:B:258:ILE:HG12	2.06	0.56
1:C:136:VAL:HG12	1:C:168:ALA:HB2	1.88	0.56
1:D:136:VAL:HG12	1:D:168:ALA:HB2	1.88	0.56
1:D:41:VAL:HG22	1:D:243:VAL:HG21	1.88	0.56
1:D:160:ILE:HD12	1:D:160:ILE:N	2.21	0.56
1:D:230:PRO:HG3	4:D:2027:HOH:O	2.05	0.56
1:A:224:TYR:CE2	1:A:228:ARG:HD3	2.41	0.56
1:B:225:ASN:HD22	1:B:228:ARG:HE	1.54	0.56
1:A:22:THR:OG1	1:A:305:VAL:HG22	2.06	0.55
1:C:97:ASP:O	1:C:119:LYS:CG	2.54	0.55
1:D:23:TYR:HE2	1:D:26:LEU:HD12	1.72	0.55
1:B:48:ASP:O	1:B:51:ARG:HG2	2.07	0.55
1:C:47:SER:OG	1:D:33:PHE:O	2.24	0.55
1:A:193:ASN:N	1:A:193:ASN:HD22	1.96	0.55
1:B:51:ARG:NH1	4:B:2014:HOH:O	2.38	0.55
1:B:125:LEU:HA	1:B:132:PRO:HD3	1.88	0.55
1:C:4:LEU:HD11	1:D:139:GLY:HA2	1.88	0.55
1:B:261:GLU:HG2	1:B:324:ASN:ND2	2.22	0.55
1:A:14:TYR:OH	1:B:57:PRO:HD2	2.07	0.55
1:D:183:SER:OG	1:D:185:VAL:HG23	2.06	0.55
1:A:116:VAL:O	1:A:116:VAL:HG13	2.07	0.55
1:C:6:LEU:C	1:C:8:ALA:H	2.10	0.55
1:A:232:ASN:H	1:A:232:ASN:HD22	1.55	0.55
1:C:48:ASP:O	1:C:51:ARG:HG2	2.06	0.55
1:B:58:ASP:HB2	1:B:238:GLN:HE21	1.71	0.55
1:B:193:ASN:ND2	1:B:193:ASN:H	2.02	0.55
1:C:121:LEU:CD1	1:C:123:VAL:HG22	2.37	0.55
1:C:315:ARG:HD2	4:C:2093:HOH:O	2.06	0.55
1:A:315:ARG:HD2	4:A:2097:HOH:O	2.06	0.54
1:C:96:PHE:CG	1:C:141:VAL:HB	2.41	0.54
1:C:9:LYS:N	4:C:2003:HOH:O	2.40	0.54
1:C:91:VAL:CG2	1:C:224:TYR:CE1	2.82	0.54
1:B:22:THR:CG2	1:B:305:VAL:HG22	2.36	0.54
1:C:299:HIS:HD2	4:C:2086:HOH:O	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:249:GLU:HA	1:D:252:GLU:OE2	2.07	0.54
1:A:262:ARG:HB2	1:A:318:ILE:HD11	1.90	0.54
1:D:110:ARG:NH1	1:D:114:LYS:NZ	2.54	0.54
1:C:255:THR:O	1:C:259:VAL:HG23	2.08	0.54
1:C:14:TYR:OH	1:D:57:PRO:HD2	2.08	0.54
1:A:88:ILE:HD13	1:A:214:VAL:CG2	2.35	0.54
1:C:5:ASP:O	1:C:8:ALA:HB3	2.07	0.54
1:C:296:LEU:HD22	1:C:300:LEU:HG	1.89	0.54
1:A:41:VAL:O	1:A:45:LEU:HD13	2.07	0.54
1:D:279:SER:O	1:D:280:ARG:HD3	2.08	0.54
1:A:58:ASP:HB2	1:A:238:GLN:NE2	2.22	0.54
1:B:262:ARG:HB2	1:B:318:ILE:CD1	2.37	0.54
1:A:295:ARG:HB3	1:A:334:PHE:CZ	2.39	0.53
1:C:129:LEU:HB3	1:C:153:HIS:CE1	2.43	0.53
1:C:56:SER:HB3	1:C:57:PRO:CD	2.35	0.53
1:A:9:LYS:HG3	1:B:226:ARG:NH2	2.23	0.53
1:B:311:ARG:NH2	1:D:249:GLU:HG2	2.22	0.53
1:D:87:GLU:O	1:D:91:VAL:HG13	2.08	0.53
1:B:262:ARG:HG3	1:B:318:ILE:HD11	1.89	0.53
1:C:10:ARG:HB3	1:C:10:ARG:NH1	2.23	0.53
1:C:119:LYS:HD3	1:C:119:LYS:N	2.23	0.53
1:D:7:ILE:C	1:D:9:LYS:N	2.60	0.53
1:D:81:VAL:HG22	1:D:82:GLY:N	2.23	0.53
1:C:239:MET:O	1:C:243:VAL:HG13	2.09	0.53
1:B:29:ASN:HB2	1:B:202:LYS:O	2.09	0.53
1:A:51:ARG:HD2	1:B:29:ASN:O	2.08	0.53
1:C:81:VAL:HG22	1:C:82:GLY:N	2.24	0.53
1:C:10:ARG:NH2	4:C:2004:HOH:O	2.42	0.53
1:B:62:ILE:HG22	1:B:66:LEU:CD2	2.39	0.53
1:A:262:ARG:HB2	1:A:318:ILE:CD1	2.39	0.53
1:A:136:VAL:HA	1:A:140:ASP:OD1	2.09	0.52
1:B:202:LYS:CE	2:B:1335:PMP:H4A2	2.39	0.52
1:B:6:LEU:O	1:B:6:LEU:HD12	2.09	0.52
1:B:308:ARG:HG3	1:B:310:PHE:CE2	2.44	0.52
1:C:71:THR:HG21	1:C:74:LEU:CB	2.31	0.52
1:C:315:ARG:HH22	3:C:1336:HSA:P	2.32	0.52
1:B:262:ARG:HB2	1:B:318:ILE:HD13	1.92	0.52
1:A:57:PRO:HG2	1:A:228:ARG:NH2	2.24	0.52
1:B:222:ASP:O	1:B:226:ARG:HG3	2.09	0.52
1:A:32:PRO:HB3	1:B:48:ASP:HA	1.92	0.52
1:D:110:ARG:NH1	1:D:114:LYS:HZ3	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:311:ARG:HG2	1:D:311:ARG:HH21	1.73	0.52
1:A:73:PHE:O	1:C:21:LYS:CG	2.57	0.52
1:D:126:THR:CG2	1:D:127:LYS:N	2.73	0.52
1:D:164:LEU:HG	1:D:191:TYR:CD2	2.45	0.52
1:B:232:ASN:H	1:B:232:ASN:ND2	2.01	0.52
1:A:9:LYS:CB	1:B:226:ARG:HD2	2.39	0.52
1:C:10:ARG:CZ	4:C:2004:HOH:O	2.56	0.52
1:C:136:VAL:HG13	1:C:140:ASP:OD1	2.11	0.52
1:D:202:LYS:HZ3	2:D:1335:PMP:H4A2	1.75	0.52
1:D:56:SER:OG	1:D:232:ASN:ND2	2.43	0.51
1:B:8:ALA:C	1:B:10:ARG:H	2.14	0.51
1:A:234:SER:HB2	1:B:209:GLN:HA	1.92	0.51
1:A:73:PHE:O	1:C:21:LYS:HG2	2.10	0.51
1:C:274:TYR:CZ	1:C:335:LYS:HB3	2.45	0.51
1:C:119:LYS:N	1:C:119:LYS:HE2	2.25	0.51
1:B:91:VAL:HG21	1:B:224:TYR:CE1	2.45	0.51
1:A:272:MET:HE1	1:A:331:LEU:C	2.30	0.51
1:D:4:LEU:HD13	1:D:7:ILE:HD12	1.92	0.51
1:D:193:ASN:HD22	1:D:193:ASN:N	2.02	0.51
1:D:277:THR:HG22	1:D:287:PHE:CE1	2.45	0.51
1:A:144:ILE:HD11	1:A:163:ILE:HD12	1.92	0.51
1:B:40:GLU:OE1	1:B:44:ARG:HD2	2.11	0.51
1:D:142:VAL:HG11	1:D:163:ILE:CG2	2.40	0.51
1:A:138:GLU:OE1	1:B:3:PRO:HD3	2.11	0.51
1:C:198:ARG:HB2	1:C:213:TYR:CZ	2.46	0.51
1:D:326:MET:O	1:D:330:GLU:HB2	2.10	0.51
1:D:43:ARG:HD2	4:D:2010:HOH:O	2.10	0.51
1:C:226:ARG:CD	1:D:8:ALA:HA	2.41	0.51
1:C:6:LEU:C	1:C:8:ALA:N	2.62	0.51
1:A:119:LYS:HD2	4:A:2043:HOH:O	2.11	0.51
1:D:157:ARG:O	1:D:161:GLU:HG3	2.10	0.51
1:C:236:VAL:HG23	1:D:209:GLN:NE2	2.26	0.50
1:D:56:SER:O	1:D:238:GLN:NE2	2.44	0.50
1:B:148:ASN:O	1:B:152:GLY:N	2.41	0.50
1:D:32:PRO:O	1:D:254:ARG:NH1	2.42	0.50
1:B:59:GLU:HB2	4:B:2020:HOH:O	2.10	0.50
1:C:226:ARG:CZ	1:D:8:ALA:HA	2.41	0.50
1:B:293:LYS:HD3	4:B:2073:HOH:O	2.10	0.50
1:C:57:PRO:CD	1:C:228:ARG:HH22	2.24	0.50
1:C:193:ASN:HB3	1:D:1:MET:HG2	1.94	0.50
1:D:57:PRO:HG2	1:D:228:ARG:HH22	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:LEU:HD21	1:B:116:VAL:CG1	2.40	0.50
1:B:261:GLU:OE2	1:B:321:ARG:HD3	2.11	0.50
1:B:174:GLU:HB3	1:B:177:TYR:HB2	1.94	0.50
1:C:142:VAL:CG1	1:C:170:VAL:HG22	2.42	0.50
1:A:91:VAL:CG2	1:A:224:TYR:HE1	2.17	0.50
1:D:91:VAL:O	1:D:95:MET:HG3	2.12	0.50
1:C:48:ASP:HA	1:D:32:PRO:HB3	1.94	0.50
1:A:110:ARG:NH2	1:A:114:LYS:HE2	2.27	0.50
1:D:59:GLU:O	1:D:63:GLU:HG3	2.12	0.50
1:D:222:ASP:O	1:D:226:ARG:HG3	2.12	0.49
1:B:180:HIS:CD2	1:B:280:ARG:HG2	2.47	0.49
1:B:24:LEU:HD22	1:B:317:THR:O	2.12	0.49
1:D:57:PRO:HG3	1:D:81:VAL:HG12	1.94	0.49
1:A:64:LYS:HB3	1:A:245:LEU:HD12	1.95	0.49
1:D:56:SER:CB	1:D:57:PRO:HD3	2.35	0.49
1:A:21:LYS:HG2	1:A:304:ASN:CB	2.41	0.49
1:C:92:MET:HE3	1:C:197:ILE:CD1	2.42	0.49
1:A:58:ASP:CB	1:A:238:GLN:NE2	2.74	0.49
1:C:119:LYS:CD	1:C:119:LYS:N	2.67	0.49
1:D:110:ARG:HH12	1:D:114:LYS:HZ1	1.60	0.49
1:B:294:GLU:HA	1:B:294:GLU:OE2	2.13	0.49
1:B:56:SER:OG	1:B:233:VAL:N	2.44	0.49
1:D:325:ASP:O	1:D:329:ARG:HG3	2.12	0.49
1:D:222:ASP:OD2	1:D:226:ARG:HD2	2.12	0.49
1:B:103:PRO:HA	1:B:104:PRO:C	2.32	0.49
1:A:24:LEU:HA	4:A:2009:HOH:O	2.11	0.49
1:D:190:LYS:HD2	1:D:190:LYS:N	2.27	0.49
1:B:61:LEU:O	1:B:65:ILE:HG12	2.13	0.49
3:C:1336:HSA:HB2	1:D:53:TYR:CZ	2.48	0.49
1:B:146:ASN:HA	1:B:147:PRO:C	2.34	0.49
1:D:131:ILE:HD12	1:D:159:GLU:HB3	1.94	0.49
1:D:106:TYR:CZ	1:D:108:CYS:HB2	2.48	0.49
1:B:65:ILE:HD13	1:B:245:LEU:HD11	1.94	0.49
1:C:234:SER:HB2	1:D:209:GLN:HA	1.95	0.48
1:D:163:ILE:O	1:D:166:THR:HG22	2.13	0.48
1:D:202:LYS:CE	2:D:1335:PMP:H4A2	2.42	0.48
1:C:60:GLU:HA	1:C:63:GLU:OE2	2.12	0.48
1:C:118:ALA:HB1	1:C:119:LYS:HE2	1.95	0.48
1:A:289:GLU:O	1:A:291:GLU:N	2.45	0.48
1:A:56:SER:O	1:A:58:ASP:N	2.46	0.48
1:A:262:ARG:O	1:A:266:LYS:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:22:THR:HG22	1:C:304:ASN:O	2.14	0.48
1:A:232:ASN:ND2	1:A:232:ASN:H	2.11	0.48
1:A:58:ASP:H	1:A:238:GLN:NE2	2.11	0.48
1:D:265:MET:O	1:D:269:LEU:HG	2.13	0.48
1:A:191:TYR:HB3	1:A:193:ASN:ND2	2.28	0.48
1:B:190:LYS:H	1:B:190:LYS:HD2	1.77	0.48
1:C:57:PRO:CG	1:C:228:ARG:NH2	2.77	0.48
1:C:98:ARG:NH2	1:C:140:ASP:OD1	2.39	0.48
1:A:12:TYR:H	1:B:226:ARG:HH21	1.62	0.48
1:C:326:MET:O	1:C:330:GLU:HG2	2.14	0.48
1:B:156:GLU:HB2	1:B:159:GLU:HG3	1.96	0.48
1:B:4:LEU:N	1:B:4:LEU:HD12	2.29	0.48
1:B:311:ARG:HH22	1:D:249:GLU:CB	2.27	0.48
1:A:261:GLU:OE1	1:A:321:ARG:HA	2.13	0.48
1:B:202:LYS:HE2	2:B:1335:PMP:H4A2	1.94	0.48
1:D:81:VAL:HG22	1:D:233:VAL:HG21	1.96	0.48
1:B:96:PHE:CD2	1:B:141:VAL:HG22	2.49	0.48
1:A:261:GLU:CG	1:A:324:ASN:ND2	2.77	0.48
1:B:320:LYS:HD2	4:B:2079:HOH:O	2.14	0.48
1:B:148:ASN:O	1:B:152:GLY:HA2	2.15	0.47
1:D:193:ASN:ND2	1:D:193:ASN:H	2.02	0.47
1:A:138:GLU:OE1	1:B:3:PRO:CD	2.62	0.47
1:D:103:PRO:HA	1:D:104:PRO:C	2.34	0.47
1:C:27:ASN:HD21	1:C:315:ARG:HD3	1.78	0.47
1:B:30:PRO:HB3	1:B:319:GLY:HA2	1.95	0.47
1:B:298:GLU:N	1:D:35:GLU:HG2	2.29	0.47
4:A:2065:HOH:O	1:B:1:MET:HG3	2.13	0.47
1:B:58:ASP:O	1:B:62:ILE:HG12	2.13	0.47
1:D:22:THR:HG23	1:D:305:VAL:HA	1.95	0.47
1:B:262:ARG:HD3	1:B:281:GLY:O	2.14	0.47
1:B:20:ASP:OD2	1:B:304:ASN:HA	2.14	0.47
1:C:188:LEU:CD1	1:C:215:VAL:HG12	2.45	0.47
1:D:233:VAL:HG12	1:D:238:GLN:HG3	1.96	0.47
1:C:142:VAL:HG13	1:C:170:VAL:HG22	1.97	0.47
1:D:259:VAL:HG23	1:D:260:GLU:N	2.30	0.47
1:B:56:SER:O	1:B:238:GLN:NE2	2.48	0.47
1:D:293:LYS:O	1:D:297:LEU:HD13	2.15	0.47
1:B:111:ILE:N	1:B:111:ILE:CD1	2.78	0.47
1:A:30:PRO:HD3	4:A:2098:HOH:O	2.14	0.47
1:B:262:ARG:CB	1:B:318:ILE:HD11	2.45	0.47
1:C:262:ARG:O	1:C:266:LYS:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:PRO:HA	1:D:173:ASP:HB3	1.96	0.46
3:C:1336:HSA:HE1	4:C:2054:HOH:O	2.15	0.46
1:B:8:ALA:O	1:B:9:LYS:HB2	2.15	0.46
1:D:2:ASN:O	1:D:3:PRO:C	2.54	0.46
1:A:296:LEU:HD22	1:A:300:LEU:CD1	2.45	0.46
1:B:290:LYS:O	1:B:294:GLU:HG2	2.14	0.46
1:C:295:ARG:HB2	1:C:334:PHE:HZ	1.80	0.46
1:B:111:ILE:N	1:B:111:ILE:HD12	2.30	0.46
1:A:41:VAL:HG23	1:A:243:VAL:HG21	1.98	0.46
1:A:256:LYS:O	1:A:259:VAL:HG22	2.16	0.46
1:C:39:ASP:CG	1:C:43:ARG:HH22	2.19	0.46
1:A:81:VAL:HG23	1:A:212:GLY:O	2.15	0.46
1:B:27:ASN:HD21	1:B:315:ARG:HD3	1.80	0.46
1:B:308:ARG:CG	1:B:310:PHE:CE2	2.99	0.46
1:C:64:LYS:HE3	1:C:246:ASP:OD1	2.15	0.46
1:D:162:ARG:NH1	1:D:165:LYS:NZ	2.64	0.46
1:B:105:THR:HA	1:B:149:ASN:O	2.16	0.46
1:A:56:SER:O	1:A:238:GLN:NE2	2.49	0.46
1:B:325:ASP:O	1:B:329:ARG:HD3	2.16	0.46
1:B:60:GLU:HG3	1:B:242:LYS:CD	2.46	0.46
1:A:226:ARG:CZ	1:B:13:PRO:HA	2.46	0.46
1:D:259:VAL:O	1:D:263:GLU:HG2	2.15	0.46
1:C:248:ARG:O	1:C:252:GLU:HG3	2.16	0.46
1:B:92:MET:HG2	1:B:95:MET:HE3	1.98	0.46
1:B:142:VAL:HG11	1:B:163:ILE:HG21	1.98	0.46
1:A:102:PHE:HA	1:A:103:PRO:HD2	1.89	0.46
1:C:44:ARG:NH2	1:C:243:VAL:HG12	2.32	0.45
1:C:270:ARG:NH1	1:C:276:ILE:HD11	2.32	0.45
1:A:321:ARG:HE	1:A:321:ARG:HB2	1.54	0.45
1:D:2:ASN:ND2	1:D:3:PRO:HD3	2.31	0.45
1:C:21:LYS:H	1:C:21:LYS:CD	2.29	0.45
1:A:332:GLU:C	1:A:334:PHE:H	2.19	0.45
1:D:4:LEU:HA	1:D:4:LEU:HD22	1.59	0.45
1:B:7:ILE:O	1:B:10:ARG:HG2	2.16	0.45
1:A:103:PRO:HA	1:A:104:PRO:C	2.37	0.45
1:D:123:VAL:HG12	1:D:132:PRO:HG3	1.98	0.45
1:D:186:ASP:HB2	4:D:2051:HOH:O	2.15	0.45
1:B:200:PHE:O	1:B:204:PHE:HB2	2.16	0.45
1:D:48:ASP:O	1:D:51:ARG:HG2	2.17	0.45
1:C:103:PRO:HA	1:C:105:THR:HG22	1.98	0.45
1:C:10:ARG:HB3	1:C:10:ARG:CZ	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:274:TYR:OH	1:C:335:LYS:HB3	2.16	0.45
1:C:145:PRO:HG2	1:C:148:ASN:OD1	2.16	0.45
1:D:110:ARG:HH12	1:D:114:LYS:NZ	2.14	0.45
1:B:324:ASN:HA	1:B:327:ILE:HD12	1.99	0.45
1:C:51:ARG:HD2	1:D:29:ASN:O	2.17	0.45
1:C:297:LEU:HA	1:C:297:LEU:HD12	1.83	0.45
1:A:88:ILE:HG21	1:A:214:VAL:HG22	1.98	0.45
1:A:145:PRO:HA	1:A:173:ASP:O	2.16	0.45
1:C:125:LEU:HA	1:C:132:PRO:HD3	1.99	0.45
1:D:52:ILE:HD12	1:D:52:ILE:C	2.38	0.45
1:C:65:ILE:O	1:C:69:LEU:CD2	2.64	0.45
1:A:73:PHE:HB2	1:C:21:LYS:HG2	1.99	0.45
1:C:295:ARG:CB	1:C:334:PHE:HZ	2.30	0.45
1:D:160:ILE:O	1:D:164:LEU:HB2	2.17	0.45
2:D:1335:PMP:H4A1	2:D:1335:PMP:H5A1	1.73	0.45
1:B:293:LYS:O	1:B:297:LEU:HD22	2.17	0.45
1:C:13:PRO:HG3	1:D:226:ARG:NH2	2.32	0.44
1:C:172:LEU:HD23	1:C:184:TYR:HB2	1.99	0.44
1:D:7:ILE:O	1:D:8:ALA:CB	2.63	0.44
1:D:256:LYS:O	1:D:259:VAL:HG22	2.18	0.44
1:C:22:THR:CG2	1:C:305:VAL:HG22	2.47	0.44
1:D:191:TYR:HB3	1:D:193:ASN:ND2	2.32	0.44
1:D:7:ILE:HG22	1:D:8:ALA:N	2.31	0.44
1:C:219:LYS:HB3	1:C:219:LYS:HZ3	1.82	0.44
1:D:98:ARG:HB3	1:D:140:ASP:OD2	2.17	0.44
1:B:157:ARG:HA	1:B:184:TYR:OH	2.18	0.44
1:D:9:LYS:NZ	1:D:9:LYS:HB3	2.32	0.44
1:B:188:LEU:HD13	1:B:194:LEU:HD23	2.00	0.44
1:A:57:PRO:CD	1:A:228:ARG:HH22	2.31	0.44
1:A:54:TYR:O	1:A:55:ASP:C	2.55	0.44
1:A:184:TYR:O	1:A:187:PHE:HB2	2.17	0.44
1:D:81:VAL:CG2	1:D:233:VAL:HG21	2.48	0.43
1:D:326:MET:HE2	1:D:327:ILE:HA	1.99	0.43
1:C:43:ARG:HD3	1:C:43:ARG:HA	1.58	0.43
1:B:265:MET:SD	1:B:324:ASN:HB3	2.58	0.43
1:A:250:ILE:HD12	1:A:250:ILE:H	1.83	0.43
1:B:179:PHE:HE2	1:B:202:LYS:HB3	1.83	0.43
1:C:172:LEU:HD12	1:C:172:LEU:HA	1.90	0.43
1:D:105:THR:HA	1:D:149:ASN:O	2.18	0.43
1:B:128:ASP:O	1:B:129:LEU:HB2	2.18	0.43
1:A:56:SER:OG	1:A:232:ASN:ND2	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:SER:HB2	1:A:198:ARG:HH21	1.84	0.43
1:A:60:GLU:CD	1:A:242:LYS:HD2	2.38	0.43
1:C:292:GLU:HG3	1:C:293:LYS:N	2.34	0.43
1:B:87:GLU:O	1:B:91:VAL:HG13	2.19	0.43
1:D:91:VAL:CG2	1:D:224:TYR:CE1	2.99	0.43
1:C:249:GLU:OE2	1:C:250:ILE:HD12	2.18	0.43
1:C:50:LEU:HD23	1:C:235:TYR:HD2	1.83	0.43
1:C:210:ARG:CD	1:D:231:PHE:O	2.65	0.43
1:C:43:ARG:CG	1:C:43:ARG:NH2	2.71	0.43
1:A:202:LYS:CE	2:A:1335:PMP:H4A2	2.47	0.43
1:B:116:VAL:O	1:B:116:VAL:CG1	2.66	0.43
1:D:276:ILE:C	1:D:276:ILE:HD13	2.39	0.43
1:C:81:VAL:HG22	1:C:82:GLY:H	1.82	0.43
1:C:116:VAL:HG12	1:C:118:ALA:HB2	2.00	0.43
1:C:257:PHE:O	1:C:261:GLU:HB2	2.18	0.43
1:A:40:GLU:OE1	1:A:247:HIS:CE1	2.71	0.43
1:B:234:SER:O	1:B:238:GLN:HG3	2.17	0.43
1:C:21:LYS:N	1:C:21:LYS:HD2	2.33	0.43
1:B:162:ARG:O	1:B:166:THR:HG23	2.19	0.43
1:D:283:PHE:HA	1:D:318:ILE:HG12	2.01	0.43
1:D:172:LEU:HD21	1:D:184:TYR:CD2	2.54	0.43
1:D:166:THR:CG2	1:D:168:ALA:H	2.32	0.43
1:A:65:ILE:HD13	1:A:245:LEU:HD11	1.99	0.43
1:C:164:LEU:HA	1:C:164:LEU:HD12	1.92	0.43
1:C:28:GLU:OE2	1:D:52:ILE:HB	2.19	0.43
1:C:320:LYS:HE3	4:C:2018:HOH:O	2.18	0.43
1:C:30:PRO:HB3	1:C:319:GLY:HA2	1.99	0.43
1:C:232:ASN:ND2	1:C:232:ASN:N	2.50	0.43
1:A:106:TYR:CZ	1:A:108:CYS:HB2	2.53	0.43
1:A:143:PHE:HA	1:A:171:ALA:HB3	2.01	0.43
1:D:126:THR:HG22	1:D:127:LYS:N	2.34	0.42
1:D:57:PRO:CG	1:D:228:ARG:NH2	2.78	0.42
1:A:272:MET:HE1	1:A:332:GLU:N	2.33	0.42
1:C:226:ARG:HG2	1:D:12:TYR:O	2.19	0.42
1:B:145:PRO:HG2	1:B:148:ASN:OD1	2.19	0.42
2:B:1335:PMP:H5A1	2:B:1335:PMP:H4A1	1.80	0.42
1:A:34:PRO:HB2	1:A:37:LEU:HG	2.01	0.42
1:B:56:SER:HB3	1:B:57:PRO:CD	2.48	0.42
1:C:27:ASN:HA	1:C:317:THR:CG2	2.49	0.42
1:A:265:MET:O	1:A:269:LEU:HG	2.18	0.42
1:B:45:LEU:HD22	1:B:45:LEU:HA	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:288:MET:HE3	1:C:293:LYS:HG3	2.02	0.42
1:D:41:VAL:CG2	1:D:243:VAL:HG21	2.49	0.42
1:A:34:PRO:O	1:A:38:VAL:HG23	2.18	0.42
1:A:112:PHE:O	1:A:116:VAL:HB	2.18	0.42
1:B:311:ARG:CZ	1:D:249:GLU:HG2	2.50	0.42
1:D:333:VAL:HG22	1:D:333:VAL:O	2.18	0.42
1:B:269:LEU:HD21	1:B:331:LEU:HD12	2.01	0.42
1:C:261:GLU:CG	1:C:324:ASN:ND2	2.74	0.42
1:D:4:LEU:HB3	4:D:2001:HOH:O	2.19	0.42
1:D:136:VAL:HG12	1:D:137:GLY:N	2.34	0.42
1:A:69:LEU:CD1	1:A:69:LEU:N	2.82	0.42
1:C:219:LYS:HB3	1:C:219:LYS:HZ2	1.84	0.42
1:A:154:VAL:HG21	1:A:180:HIS:NE2	2.35	0.42
1:C:261:GLU:HG2	1:C:324:ASN:HD22	1.80	0.42
1:D:288:MET:HB2	1:D:293:LYS:HG3	2.01	0.42
1:D:261:GLU:OE2	1:D:321:ARG:HA	2.19	0.42
1:C:65:ILE:HG22	1:C:69:LEU:HD23	2.02	0.42
1:A:297:LEU:O	1:A:301:ARG:HB2	2.20	0.42
1:D:268:ALA:O	1:D:272:MET:HG3	2.19	0.42
1:D:58:ASP:HB2	1:D:238:GLN:HE21	1.73	0.42
1:C:95:MET:O	1:D:4:LEU:CD2	2.64	0.42
1:C:123:VAL:HG21	1:C:134:VAL:HG11	2.02	0.42
1:A:252:GLU:HG2	4:A:2036:HOH:O	2.19	0.42
1:D:56:SER:O	1:D:238:GLN:CD	2.58	0.42
1:B:94:LEU:CD2	1:B:116:VAL:CG1	2.91	0.42
1:C:138:GLU:OE1	1:D:3:PRO:HD2	2.20	0.42
1:C:179:PHE:CD1	1:C:255:THR:HG23	2.55	0.42
1:D:56:SER:CB	1:D:232:ASN:ND2	2.83	0.42
1:C:57:PRO:CD	1:C:228:ARG:NH2	2.83	0.42
1:D:74:LEU:HD21	4:D:2052:HOH:O	2.20	0.42
1:C:197:ILE:HD13	1:C:214:VAL:HG13	2.01	0.42
1:D:106:TYR:CE2	1:D:108:CYS:HB2	2.55	0.42
1:A:164:LEU:HD13	1:A:170:VAL:CG2	2.49	0.42
1:A:50:LEU:HA	1:B:208:ALA:HB3	2.02	0.42
1:C:33:PHE:CD1	1:C:206:LEU:HD13	2.55	0.42
1:D:88:ILE:O	1:D:92:MET:HG3	2.20	0.41
1:C:119:LYS:HG2	1:C:119:LYS:O	2.20	0.41
1:C:78:ASN:O	1:C:215:VAL:HA	2.20	0.41
1:C:308:ARG:HG2	1:C:310:PHE:CE2	2.54	0.41
1:D:141:VAL:HA	1:D:169:PHE:O	2.20	0.41
1:D:116:VAL:HG13	1:D:116:VAL:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:160:ILE:CD1	1:D:160:ILE:N	2.82	0.41
1:B:21:LYS:HB2	1:B:304:ASN:HB3	2.02	0.41
1:D:2:ASN:HD22	1:D:3:PRO:HD3	1.84	0.41
1:C:79:VAL:HA	1:C:214:VAL:O	2.21	0.41
1:B:210:ARG:NH2	2:B:1335:PMP:O1P	2.48	0.41
1:D:81:VAL:HG23	1:D:212:GLY:O	2.20	0.41
1:A:57:PRO:CD	1:A:228:ARG:NH2	2.84	0.41
1:A:254:ARG:O	1:A:258:ILE:HG12	2.21	0.41
1:C:210:ARG:NH2	2:C:1335:PMP:O2P	2.37	0.41
1:D:326:MET:O	1:D:326:MET:CE	2.69	0.41
1:B:45:LEU:HD13	1:B:235:TYR:HE2	1.86	0.41
1:A:56:SER:OG	1:A:233:VAL:HB	2.20	0.41
1:B:184:TYR:O	1:B:187:PHE:HB2	2.20	0.41
1:D:331:LEU:HD23	1:D:331:LEU:HA	1.93	0.41
1:C:56:SER:OG	1:C:232:ASN:ND2	2.54	0.41
1:B:58:ASP:N	1:B:238:GLN:HE22	2.19	0.41
1:A:79:VAL:HA	1:A:214:VAL:O	2.21	0.41
1:A:263:GLU:OE1	1:A:266:LYS:HE2	2.20	0.41
1:C:297:LEU:O	1:C:301:ARG:HG3	2.20	0.41
1:C:335:LYS:HG3	1:C:335:LYS:OXT	2.20	0.41
1:D:146:ASN:HA	1:D:147:PRO:C	2.41	0.41
1:D:56:SER:HB3	1:D:57:PRO:CD	2.38	0.41
1:C:136:VAL:CG1	1:C:137:GLY:N	2.82	0.41
1:A:267:SER:HA	1:A:270:ARG:NH1	2.36	0.41
1:B:148:ASN:O	1:B:152:GLY:CA	2.69	0.41
1:C:61:LEU:O	1:C:65:ILE:HG12	2.20	0.41
1:B:7:ILE:C	1:B:8:ALA:O	2.58	0.41
1:D:296:LEU:HD11	1:D:300:LEU:HD11	2.01	0.41
1:C:53:TYR:CZ	3:D:1336:HSA:HB2	2.55	0.41
1:B:198:ARG:HB2	1:B:213:TYR:CZ	2.56	0.41
1:D:179:PHE:CD1	1:D:255:THR:HG23	2.56	0.41
1:A:308:ARG:HA	1:A:308:ARG:HD3	1.74	0.41
1:A:250:ILE:N	1:A:250:ILE:HD12	2.35	0.41
1:D:22:THR:CG2	1:D:305:VAL:HG22	2.50	0.41
1:C:96:PHE:CD2	1:C:141:VAL:HB	2.56	0.41
1:C:299:HIS:CD2	4:C:2086:HOH:O	2.71	0.41
1:D:73:PHE:CD2	1:D:189:LYS:HE3	2.56	0.41
1:A:30:PRO:CG	1:A:258:ILE:HD11	2.51	0.40
1:D:20:ASP:OD1	1:D:304:ASN:HB3	2.20	0.40
1:B:110:ARG:O	1:B:114:LYS:HG3	2.20	0.40
1:D:54:TYR:O	1:D:55:ASP:C	2.59	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:232:ASN:N	1:D:232:ASN:ND2	2.64	0.40
1:A:263:GLU:OE1	1:A:263:GLU:HA	2.21	0.40
1:C:199:THR:HA	1:C:212:GLY:HA2	2.03	0.40
1:D:126:THR:C	1:D:128:ASP:N	2.73	0.40
1:B:58:ASP:H	1:B:238:GLN:NE2	2.20	0.40
1:C:326:MET:HA	1:C:329:ARG:CZ	2.51	0.40
1:B:190:LYS:HE3	4:B:2044:HOH:O	2.21	0.40
1:D:262:ARG:O	1:D:266:LYS:HG3	2.21	0.40
1:B:66:LEU:HD13	1:B:66:LEU:HA	1.91	0.40
1:A:57:PRO:CG	1:A:228:ARG:NH2	2.85	0.40
1:B:202:LYS:HD2	1:B:202:LYS:N	2.36	0.40
1:B:22:THR:HG22	1:B:304:ASN:O	2.21	0.40
1:C:300:LEU:HA	1:C:300:LEU:HD23	1.90	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	328/335 (98%)	308 (94%)	18 (6%)	2 (1%)	30	40
1	B	333/335 (99%)	319 (96%)	12 (4%)	2 (1%)	30	40
1	C	333/335 (99%)	313 (94%)	18 (5%)	2 (1%)	30	40
1	D	333/335 (99%)	309 (93%)	19 (6%)	5 (2%)	13	15
All	All	1327/1340 (99%)	1249 (94%)	67 (5%)	11 (1%)	24	33

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	290	LYS
1	C	4	LEU

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Mol	Chain	Res	Type
1	C	10	ARG
1	D	2	ASN
1	D	3	PRO
1	D	4	LEU
1	B	290	LYS
1	A	11	ALA
1	B	334	PHE
1	D	10	ARG
1	D	210	ARG

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	295/301 (98%)	273 (92%)	22 (8%)	17	24
1	B	300/301 (100%)	279 (93%)	21 (7%)	19	27
1	C	300/301 (100%)	280 (93%)	20 (7%)	20	29
1	D	300/301 (100%)	274 (91%)	26 (9%)	13	17
All	All	1195/1204 (99%)	1106 (93%)	89 (7%)	17	24

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ILE
1	A	9	LYS
1	A	12	TYR
1	A	14	TYR
1	A	18	LYS
1	A	20	ASP
1	A	36	ASP
1	A	43	ARG
1	A	45	LEU
1	A	66	LEU
1	A	101	PHE
1	A	110	ARG

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Mol	Chain	Res	Type
1	A	116	VAL
1	A	150	PRO
1	A	164	LEU
1	A	166	THR
1	A	193	ASN
1	A	202	LYS
1	A	206	LEU
1	A	232	ASN
1	A	296	LEU
1	A	311	ARG
1	B	6	LEU
1	B	22	THR
1	B	45	LEU
1	B	66	LEU
1	B	69	LEU
1	B	101	PHE
1	B	135	ASN
1	B	141	VAL
1	B	150	PRO
1	B	186	ASP
1	B	187	PHE
1	B	188	LEU
1	B	190	LYS
1	B	193	ASN
1	B	206	LEU
1	B	214	VAL
1	B	232	ASN
1	B	280	ARG
1	B	295	ARG
1	B	297	LEU
1	B	326	MET
1	C	6	LEU
1	C	10	ARG
1	C	43	ARG
1	C	102	PHE
1	C	116	VAL
1	C	119	LYS
1	C	123	VAL
1	C	150	PRO
1	C	161	GLU
1	C	166	THR
1	C	193	ASN

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Mol	Chain	Res	Type
1	C	202	LYS
1	C	206	LEU
1	C	232	ASN
1	C	249	GLU
1	C	289	GLU
1	C	290	LYS
1	C	296	LEU
1	C	297	LEU
1	C	333	VAL
1	D	2	ASN
1	D	3	PRO
1	D	4	LEU
1	D	10	ARG
1	D	22	THR
1	D	66	LEU
1	D	69	LEU
1	D	101	PHE
1	D	116	VAL
1	D	121	LEU
1	D	127	LYS
1	D	130	ARG
1	D	150	PRO
1	D	158	GLU
1	D	166	THR
1	D	186	ASP
1	D	188	LEU
1	D	193	ASN
1	D	214	VAL
1	D	232	ASN
1	D	249	GLU
1	D	250	ILE
1	D	275	ARG
1	D	276	ILE
1	D	290	LYS
1	D	326	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	77	ASN
1	A	193	ASN

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Mol	Chain	Res	Type
1	A	209	GLN
1	A	225	ASN
1	A	232	ASN
1	A	238	GLN
1	A	247	HIS
1	B	27	ASN
1	B	180	HIS
1	B	193	ASN
1	B	209	GLN
1	B	225	ASN
1	B	232	ASN
1	B	238	GLN
1	C	27	ASN
1	C	77	ASN
1	C	135	ASN
1	C	193	ASN
1	C	209	GLN
1	C	225	ASN
1	C	232	ASN
1	C	238	GLN
1	C	299	HIS
1	C	304	ASN
1	D	27	ASN
1	D	193	ASN
1	D	209	GLN
1	D	232	ASN
1	D	238	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PMP	A	1335	3	16,16,16	1.33	1 (6%)	20,23,23	1.90	3 (15%)
3	HSA	A	1336	2	9,13,14	2.32	3 (33%)	12,17,19	1.81	3 (25%)
2	PMP	B	1335	3	16,16,16	1.35	1 (6%)	20,23,23	1.97	4 (20%)
3	HSA	B	1336	2	9,13,14	2.33	3 (33%)	12,17,19	1.81	3 (25%)
2	PMP	C	1335	3	16,16,16	1.33	1 (6%)	20,23,23	2.02	3 (15%)
3	HSA	C	1336	2	9,13,14	2.34	3 (33%)	12,17,19	1.81	3 (25%)
2	PMP	D	1335	3	16,16,16	1.35	1 (6%)	20,23,23	1.88	3 (15%)
3	HSA	D	1336	2	9,13,14	2.32	3 (33%)	12,17,19	1.79	3 (25%)

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
2	PMP	A	1335	3	-	0/8/8/8	0/1/1/1
3	HSA	A	1336	2	1/1/1/2	0/7/8/10	0/1/1/1
2	PMP	B	1335	3	-	0/8/8/8	0/1/1/1
3	HSA	B	1336	2	1/1/1/2	0/7/8/10	0/1/1/1
2	PMP	C	1335	3	-	0/8/8/8	0/1/1/1
3	HSA	C	1336	2	1/1/1/2	0/7/8/10	0/1/1/1
2	PMP	D	1335	3	-	0/8/8/8	0/1/1/1
3	HSA	D	1336	2	1/1/1/2	0/7/8/10	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1336	HSA	CB-CG	-4.96	1.39	1.51
3	C	1336	HSA	CB-CG	-4.96	1.39	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1336	HSA	CB-CG	-4.95	1.39	1.51
3	D	1336	HSA	CB-CG	-4.89	1.39	1.51
3	D	1336	HSA	CA-CB	-2.58	1.39	1.52
3	B	1336	HSA	CA-CB	-2.58	1.39	1.52
3	C	1336	HSA	CA-CB	-2.56	1.39	1.52
3	A	1336	HSA	CA-CB	-2.56	1.39	1.52
2	C	1335	PMP	P-O1P	3.04	1.61	1.51
2	D	1335	PMP	P-O1P	3.07	1.61	1.51
2	A	1335	PMP	P-O1P	3.08	1.61	1.51
2	B	1335	PMP	P-O1P	3.11	1.61	1.51
3	B	1336	HSA	P-OP1	3.13	1.61	1.51
3	A	1336	HSA	P-OP1	3.15	1.61	1.51
3	C	1336	HSA	P-OP1	3.21	1.61	1.51
3	D	1336	HSA	P-OP1	3.22	1.61	1.51

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1336	HSA	CB-CG-CD2	-4.47	122.63	129.55
3	D	1336	HSA	CB-CG-CD2	-4.34	122.84	129.55
3	C	1336	HSA	CB-CG-CD2	-4.31	122.89	129.55
3	A	1336	HSA	CB-CG-CD2	-4.30	122.91	129.55
2	B	1335	PMP	C5-C6-N1	-2.05	120.29	123.86
3	B	1336	HSA	OP3-P-OP4	2.16	112.78	106.56
3	C	1336	HSA	OP3-P-OP4	2.17	112.82	106.56
3	D	1336	HSA	OP3-P-OP4	2.20	112.91	106.56
3	A	1336	HSA	OP3-P-OP4	2.48	113.71	106.56
2	C	1335	PMP	O3P-P-O4P	2.56	113.95	106.56
3	B	1336	HSA	CB-CA-C	2.59	122.44	113.29
2	C	1335	PMP	C6-C5-C4	2.64	120.06	118.09
2	D	1335	PMP	O3P-P-O4P	2.64	114.18	106.56
3	A	1336	HSA	CB-CA-C	2.66	122.71	113.29
2	B	1335	PMP	O3P-P-O4P	2.75	114.47	106.56
3	D	1336	HSA	CB-CA-C	2.76	123.06	113.29
3	C	1336	HSA	CB-CA-C	2.80	123.19	113.29
2	A	1335	PMP	O3P-P-O4P	2.94	115.03	106.56
2	A	1335	PMP	C6-C5-C4	3.04	120.36	118.09
2	D	1335	PMP	C6-C5-C4	3.45	120.66	118.09
2	B	1335	PMP	C6-C5-C4	3.63	120.80	118.09
2	D	1335	PMP	O4P-C5A-C5	5.66	118.35	108.99
2	B	1335	PMP	O4P-C5A-C5	5.92	118.77	108.99
2	A	1335	PMP	O4P-C5A-C5	6.14	119.15	108.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1335	PMP	O4P-C5A-C5	7.23	120.95	108.99

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	1336	HSA	CA
3	B	1336	HSA	CA
3	C	1336	HSA	CA
3	D	1336	HSA	CA

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1335	PMP	4	0
3	A	1336	HSA	1	0
2	B	1335	PMP	4	0
2	C	1335	PMP	3	0
3	C	1336	HSA	3	0
2	D	1335	PMP	5	0
3	D	1336	HSA	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	330/335 (98%)	0.27	19 (5%) 26 31	13, 25, 60, 99	0
1	B	335/335 (100%)	0.11	9 (2%) 58 61	14, 26, 48, 65	1 (0%)
1	C	335/335 (100%)	0.42	19 (5%) 27 31	15, 31, 56, 83	0
1	D	335/335 (100%)	0.39	24 (7%) 18 21	18, 31, 61, 78	1 (0%)
All	All	1335/1340 (99%)	0.30	71 (5%) 30 34	13, 29, 56, 99	2 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	ALA	13.2
1	A	16	THR	7.9
1	C	1	MET	7.5
1	A	12	TYR	7.2
1	C	7	ILE	6.3
1	A	13	PRO	6.1
1	C	2	ASN	5.5
1	D	4	LEU	5.2
1	A	10	ARG	4.8
1	D	333	VAL	4.7
1	C	56	SER	4.6
1	A	18	LYS	4.6
1	D	302	THR	4.6
1	A	14	TYR	4.5
1	B	333	VAL	4.4
1	C	3	PRO	4.1
1	C	8	ALA	4.1
1	C	16	THR	4.1
1	D	271	GLU	4.0
1	D	18	LYS	4.0
1	B	2	ASN	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	13	PRO	3.7
1	A	6	LEU	3.7
1	D	2	ASN	3.6
1	A	7	ILE	3.6
1	C	334	PHE	3.5
1	D	1	MET	3.5
1	D	12	TYR	3.4
1	C	18	LYS	3.4
1	A	15	GLU	3.3
1	A	17	GLU	3.2
1	D	56	SER	3.1
1	D	316	ILE	3.0
1	C	335	LYS	3.0
1	D	8	ALA	3.0
1	B	181	GLY	2.9
1	B	5	ASP	2.8
1	D	290	LYS	2.8
1	D	6	LEU	2.8
1	C	70	ASP	2.6
1	C	331	LEU	2.5
1	B	182	GLU	2.5
1	D	17	GLU	2.5
1	A	9	LYS	2.5
1	D	21	LYS	2.4
1	B	127	LYS	2.4
1	A	301	ARG	2.4
1	A	291	GLU	2.4
1	C	4	LEU	2.3
1	C	135	ASN	2.3
1	A	329	ARG	2.3
1	D	273	GLY	2.3
1	C	294	GLU	2.3
1	C	17	GLU	2.2
1	C	333	VAL	2.2
1	C	15	GLU	2.2
1	D	291	GLU	2.2
1	B	1	MET	2.2
1	D	158	GLU	2.2
1	B	138	GLU	2.1
1	A	290	LYS	2.1
1	C	19	ARG	2.1
1	D	289	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	157	ARG	2.1
1	D	133	GLU	2.1
1	B	128	ASP	2.1
1	D	335	LYS	2.1
1	A	56	SER	2.0
1	A	318	ILE	2.0
1	D	3	PRO	2.0
1	A	170	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	HSA	B	1336	13/14	0.90	0.17	2.35	37,48,58,59	0
3	HSA	A	1336	13/14	0.92	0.16	1.56	31,38,50,52	1
2	PMP	B	1335	16/16	0.97	0.14	0.87	16,40,47,50	0
3	HSA	D	1336	13/14	0.92	0.15	0.71	39,43,48,52	2
3	HSA	C	1336	13/14	0.92	0.15	0.64	34,41,51,51	0
2	PMP	D	1335	16/16	0.97	0.12	-0.04	20,25,29,37	0
2	PMP	C	1335	16/16	0.96	0.11	-0.25	15,27,36,39	0
2	PMP	A	1335	16/16	0.98	0.11	-0.61	17,28,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.