



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 02:16 PM BST

PDB ID : 6VTG
Title : Naegleria gruberi RNA ligase E227A mutant apo
Authors : Unciuleac, M.C.; Goldgur, Y.; Shuman, S.
Deposited on : 2020-02-12
Resolution : 2.49 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

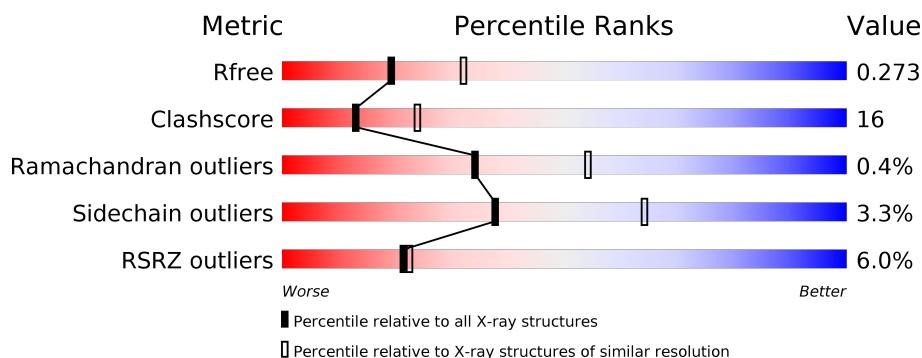
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	340	<div> <div>6%</div> <div>64%</div> <div>34%</div> <div>.</div> </div>
1	B	340	<div> <div>3%</div> <div>68%</div> <div>29%</div> <div>..</div> </div>
1	C	340	<div> <div>3%</div> <div>79%</div> <div>19%</div> <div>..</div> </div>
1	D	340	<div> <div>11%</div> <div>61%</div> <div>35%</div> <div>..</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10944 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 RNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2713	1708	480	515	10			
1	B	340	Total	C	N	O	S	0	0	0
			2713	1708	480	515	10			
1	C	340	Total	C	N	O	S	0	0	0
			2713	1708	480	515	10			
1	D	340	Total	C	N	O	S	0	0	0
			2713	1708	480	515	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP D2W2Z5
A	227	ALA	GLU	engineered mutation	UNP D2W2Z5
B	0	SER	-	expression tag	UNP D2W2Z5
B	227	ALA	GLU	engineered mutation	UNP D2W2Z5
C	0	SER	-	expression tag	UNP D2W2Z5
C	227	ALA	GLU	engineered mutation	UNP D2W2Z5
D	0	SER	-	expression tag	UNP D2W2Z5
D	227	ALA	GLU	engineered mutation	UNP D2W2Z5

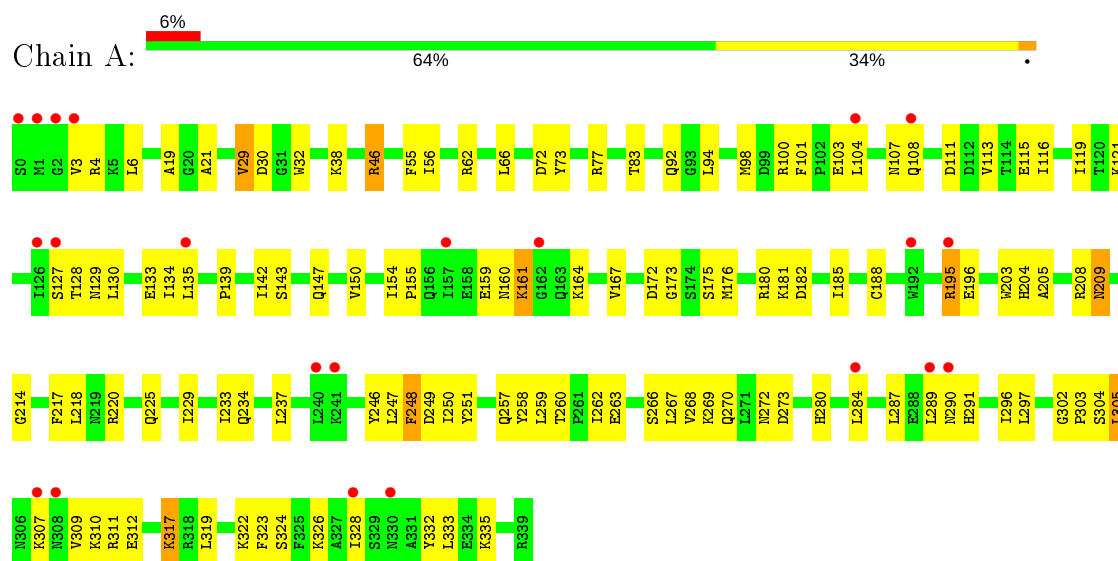
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	O	0	0
			5	5		
2	B	45	Total	O	0	0
			45	45		
2	C	35	Total	O	0	0
			35	35		
2	D	7	Total	O	0	0
			7	7		

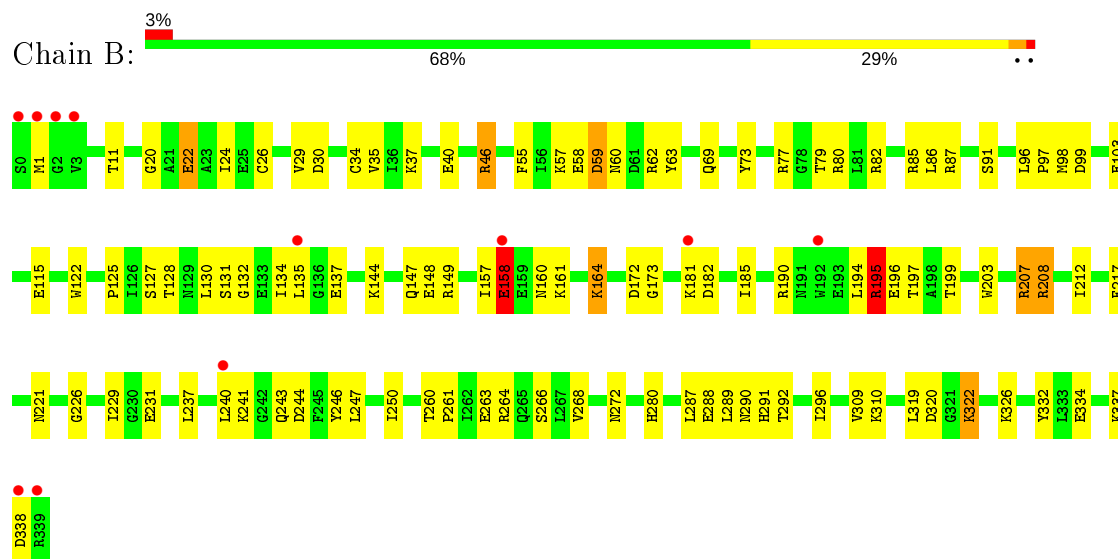
3 Residue-property plots [i](#)

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.

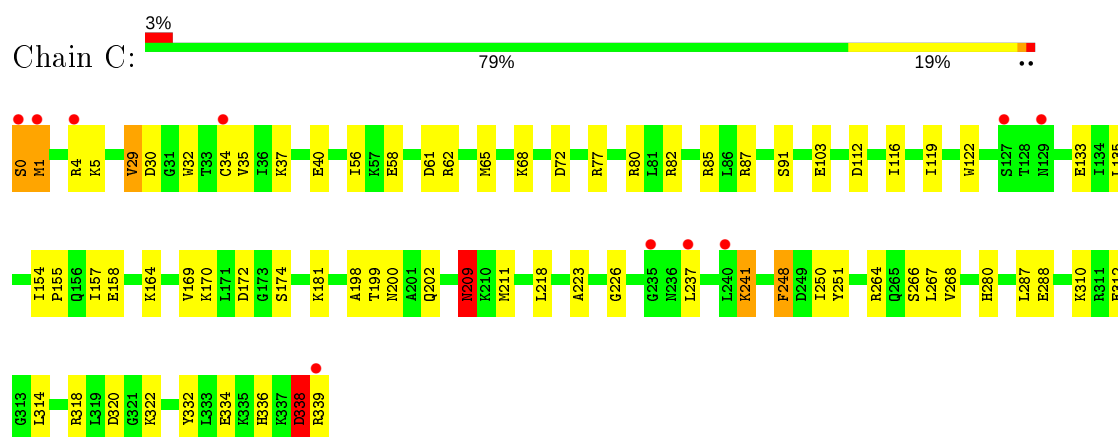
• Molecule 1: RNA ligase



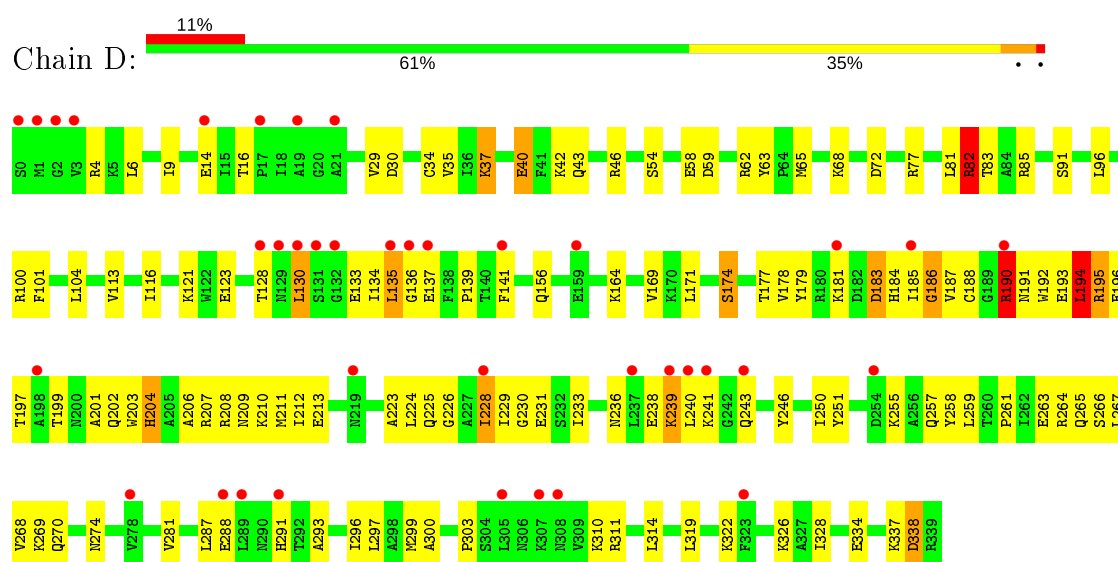
• Molecule 1: RNA ligase



• Molecule 1: RNA ligase



• Molecule 1: RNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	124.10Å 104.04Å 119.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.73 – 2.49 29.73 – 2.49	Depositor EDS
% Data completeness (in resolution range)	97.8 (29.73-2.49) 93.3 (29.73-2.49)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.215 , 0.273 0.215 , 0.273	Depositor DCC
R_{free} test set	1983 reflections (3.70%)	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 62.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.019 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10944	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.05 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.6568e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	0.52	0/2760	0.80	6/3717 (0.2%)
1	B	0.57	1/2760 (0.0%)	0.84	11/3717 (0.3%)
1	C	0.56	0/2760	0.84	6/3717 (0.2%)
1	D	0.65	5/2760 (0.2%)	0.99	15/3717 (0.4%)
All	All	0.58	6/11040 (0.1%)	0.87	38/14868 (0.3%)

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	2
1	B	0	3
1	C	0	4
1	D	0	5
All	All	0	14

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	186	GLY	CA-C	9.29	1.66	1.51
1	D	82	ARG	CB-CG	-6.99	1.33	1.52
1	D	190	ARG	CB-CG	5.95	1.68	1.52
1	D	190	ARG	CG-CD	5.90	1.66	1.51
1	B	26	CYS	CB-SG	-5.76	1.72	1.81
1	D	190	ARG	CZ-NH1	-5.37	1.26	1.33

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	190	ARG	NE-CZ-NH1	-15.10	112.75	120.30
1	C	4	ARG	NE-CZ-NH2	14.31	127.45	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	4	ARG	NE-CZ-NH1	-13.68	113.46	120.30
1	D	190	ARG	CD-NE-CZ	11.85	140.19	123.60
1	D	82	ARG	CG-CD-NE	-10.53	89.68	111.80
1	D	82	ARG	NE-CZ-NH2	-9.74	115.43	120.30
1	A	164	LYS	CD-CE-NZ	9.62	133.84	111.70
1	B	158	GLU	CA-CB-CG	-9.40	92.73	113.40
1	D	37	LYS	CA-CB-CG	-8.81	94.01	113.40
1	B	164	LYS	CB-CG-CD	8.58	133.92	111.60
1	B	164	LYS	CD-CE-NZ	-8.40	92.39	111.70
1	D	190	ARG	CA-CB-CG	8.19	131.42	113.40
1	C	338	ASP	CB-CG-OD2	-8.04	111.06	118.30
1	D	183	ASP	CB-CG-OD1	7.72	125.25	118.30
1	D	135	LEU	CA-CB-CG	7.70	133.02	115.30
1	D	228	ILE	CG1-CB-CG2	-7.50	94.90	111.40
1	A	195	ARG	NE-CZ-NH1	6.98	123.79	120.30
1	B	322	LYS	CD-CE-NZ	6.97	127.73	111.70
1	D	186	GLY	O-C-N	6.80	133.59	122.70
1	A	135	LEU	CA-CB-CG	6.29	129.77	115.30
1	C	338	ASP	CB-CG-OD1	6.24	123.91	118.30
1	D	183	ASP	CB-CG-OD2	-6.18	112.74	118.30
1	B	310	LYS	CD-CE-NZ	-6.17	97.50	111.70
1	D	194	LEU	C-N-CA	6.11	136.97	121.70
1	B	322	LYS	CB-CG-CD	5.83	126.75	111.60
1	B	22	GLU	CA-CB-CG	5.78	126.12	113.40
1	D	190	ARG	CB-CG-CD	5.77	126.60	111.60
1	C	4	ARG	CG-CD-NE	5.76	123.91	111.80
1	A	305	LEU	CA-CB-CG	5.52	127.99	115.30
1	D	239	LYS	CD-CE-NZ	-5.41	99.25	111.70
1	D	130	LEU	CA-CB-CG	-5.38	102.93	115.30
1	B	195	ARG	CB-CG-CD	5.37	125.56	111.60
1	A	164	LYS	CA-CB-CG	5.36	125.20	113.40
1	B	322	LYS	CG-CD-CE	-5.36	95.82	111.90
1	C	209	ASN	CB-CA-C	-5.24	99.92	110.40
1	B	195	ARG	N-CA-CB	5.20	119.96	110.60
1	B	59	ASP	CB-CG-OD1	5.06	122.86	118.30
1	A	317	LYS	CA-CB-CG	5.02	124.45	113.40

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	181	LYS	Peptide

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Mol	Chain	Res	Type	Group
1	A	29	VAL	Peptide
1	B	158	GLU	Sidechain
1	B	181	LYS	Peptide
1	B	29	VAL	Peptide
1	C	0	SER	Peptide
1	C	181	LYS	Peptide
1	C	29	VAL	Peptide
1	C	338	ASP	Peptide
1	D	128	THR	Peptide
1	D	190	ARG	Sidechain
1	D	194	LEU	Peptide
1	D	238	GLU	Peptide
1	D	29	VAL	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2713	0	2713	103	2
1	B	2713	0	2713	74	0
1	C	2713	0	2713	58	0
1	D	2713	0	2713	126	2
2	A	5	0	0	6	0
2	B	45	0	0	11	0
2	C	35	0	0	6	0
2	D	7	0	0	4	0
All	All	10944	0	10852	355	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 16.

All (355) 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:4:ARG:NH2	1:D:121:LYS:HB2	1.52	1.20
1:A:269:LYS:N	2:A:401:HOH:O	1.85	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLN:HB3	1:A:111:ASP:OD1	1.56	1.06
1:B:57:LYS:NZ	2:B:401:HOH:O	1.89	1.04
1:A:108:GLN:HB3	1:A:111:ASP:CG	1.77	1.04
1:A:108:GLN:CB	1:A:111:ASP:OD1	2.17	0.93
1:A:317:LYS:HB3	1:A:324:SER:HB3	1.50	0.92
1:D:58:GLU:OE1	1:D:77:ARG:NH1	2.02	0.90
1:D:137:GLU:HG3	1:D:192:TRP:CE3	2.07	0.90
1:D:4:ARG:NH2	1:D:121:LYS:CB	2.35	0.89
1:A:257:GLN:NE2	2:A:402:HOH:O	2.05	0.88
1:C:5:LYS:HE3	1:C:112:ASP:HB2	1.52	0.88
1:A:100:ARG:NH2	2:A:403:HOH:O	2.08	0.86
1:D:288:GLU:O	1:D:288:GLU:HG2	1.75	0.85
1:D:196:GLU:OE2	2:D:401:HOH:O	1.94	0.84
1:D:4:ARG:HH21	1:D:121:LYS:HB2	1.39	0.82
1:B:1:MET:O	2:B:402:HOH:O	1.96	0.81
1:D:82:ARG:HH22	1:D:337:LYS:CB	1.96	0.79
1:C:172:ASP:OD2	1:C:332:TYR:OH	2.02	0.78
1:B:132:GLY:HA3	1:B:194:LEU:HD23	1.64	0.78
1:C:338:ASP:HB2	1:C:339:ARG:HH11	1.47	0.78
1:D:190:ARG:NH1	1:D:191:ASN:HD21	1.82	0.77
1:D:259:LEU:HD21	1:D:267:LEU:HD12	1.67	0.77
1:B:208:ARG:NH2	1:B:244:ASP:OD2	2.19	0.76
1:D:137:GLU:HG3	1:D:192:TRP:HE3	1.51	0.75
1:A:234:GLN:OE1	1:A:234:GLN:N	2.20	0.75
1:A:161:LYS:HD2	1:A:290:ASN:HA	1.67	0.74
1:C:338:ASP:HB2	1:C:339:ARG:NH1	2.02	0.74
1:A:62:ARG:NH2	1:A:103:GLU:OE2	2.21	0.74
1:A:233:ILE:HA	1:A:234:GLN:OE1	1.89	0.73
1:D:82:ARG:NH2	1:D:337:LYS:O	2.22	0.73
1:C:339:ARG:HH11	1:C:339:ARG:HG2	1.52	0.73
1:D:134:ILE:HG12	1:D:192:TRP:HB3	1.68	0.73
1:B:24:ILE:O	2:B:403:HOH:O	2.07	0.72
1:A:205:ALA:O	1:A:209:ASN:ND2	2.21	0.72
1:D:130:LEU:HD22	1:D:202:GLN:HG3	1.71	0.72
1:C:209:ASN:ND2	1:C:211:MET:SD	2.63	0.71
1:D:82:ARG:HH22	1:D:337:LYS:HB3	1.54	0.71
1:D:113:VAL:HA	1:D:116:ILE:HD13	1.71	0.71
1:A:103:GLU:HB3	1:D:310:LYS:HE3	1.73	0.71
1:D:82:ARG:HD3	1:D:83:THR:O	1.91	0.71
1:D:82:ARG:NH2	1:D:337:LYS:HG2	2.06	0.70
1:D:210:LYS:HZ1	1:D:212:ILE:HB	1.57	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:LYS:CE	1:A:335:LYS:HD2	2.22	0.70
1:D:223:ALA:HB3	1:D:251:TYR:HB3	1.74	0.70
1:A:269:LYS:CA	2:A:401:HOH:O	2.32	0.70
1:A:268:VAL:HG12	1:A:272:ASN:HD21	1.58	0.68
1:B:60:ASN:OD1	2:B:404:HOH:O	2.12	0.68
1:A:310:LYS:HE3	1:A:335:LYS:HD2	1.76	0.68
1:D:225:GLN:NE2	1:D:258:TYR:OH	2.27	0.67
1:D:190:ARG:NH1	1:D:191:ASN:ND2	2.42	0.67
1:D:14:GLU:OE2	1:D:16:THR:HG22	1.95	0.66
1:D:177:THR:HG22	1:D:225:GLN:HG2	1.77	0.66
1:B:250:ILE:HD12	1:B:264:ARG:HG3	1.78	0.66
1:B:172:ASP:OD2	1:B:332:TYR:OH	2.14	0.66
1:C:209:ASN:OD1	1:C:211:MET:N	2.29	0.66
1:B:182:ASP:OD1	1:B:221:ASN:ND2	2.21	0.66
1:D:116:ILE:H	1:D:116:ILE:HD12	1.59	0.66
1:C:62:ARG:NH2	1:C:103:GLU:OE2	2.29	0.65
1:D:259:LEU:HD23	1:D:264:ARG:HA	1.78	0.65
1:D:82:ARG:NH2	1:D:337:LYS:CB	2.59	0.65
1:D:85:ARG:NH1	1:D:334:GLU:OE2	2.29	0.65
1:D:190:ARG:NH1	1:D:191:ASN:OD1	2.30	0.65
1:D:133:GLU:O	1:D:195:ARG:HG3	1.98	0.64
1:B:125:PRO:O	2:B:405:HOH:O	2.14	0.64
1:A:269:LYS:CB	2:A:401:HOH:O	2.44	0.64
1:B:11:THR:O	2:B:406:HOH:O	2.15	0.64
1:A:3:VAL:HG12	1:A:4:ARG:H	1.63	0.63
1:D:82:ARG:HH22	1:D:337:LYS:CG	2.11	0.63
1:A:237:LEU:O	1:A:309:VAL:HG11	1.98	0.62
1:A:154:ILE:HG23	1:A:155:PRO:HD3	1.81	0.62
1:B:80:ARG:NH1	2:B:411:HOH:O	2.32	0.62
1:A:260:THR:OG1	1:A:263:GLU:HG3	1.99	0.62
1:D:263:GLU:O	1:D:267:LEU:HG	1.99	0.61
1:C:199:THR:HG23	2:C:403:HOH:O	1.99	0.61
1:D:230:GLY:O	1:D:236:ASN:HB2	2.00	0.61
1:A:167:VAL:HG23	1:A:284:LEU:HB2	1.83	0.61
1:B:148:GLU:OE1	1:B:149:ARG:N	2.31	0.61
1:B:260:THR:OG1	1:B:263:GLU:HG3	2.01	0.61
1:D:137:GLU:HG3	1:D:192:TRP:CZ3	2.36	0.60
1:B:247:LEU:HD23	1:B:280:HIS:CD2	2.36	0.60
1:D:82:ARG:HH22	1:D:337:LYS:HG2	1.65	0.60
1:D:207:ARG:HD2	2:D:401:HOH:O	2.01	0.60
1:B:20:GLY:HA3	1:B:87:ARG:HH21	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:LEU:HD23	1:B:280:HIS:HD2	1.64	0.60
1:C:87:ARG:NE	2:C:401:HOH:O	2.05	0.59
1:D:210:LYS:HE3	1:D:213:GLU:OE1	2.02	0.59
1:A:161:LYS:CD	1:A:290:ASN:HA	2.32	0.59
1:D:169:VAL:HG13	1:D:311:ARG:HD2	1.85	0.59
1:A:173:GLY:HA3	1:A:229:ILE:HG22	1.84	0.58
1:B:135:LEU:HD12	1:B:195:ARG:HG3	1.84	0.58
1:A:56:ILE:HA	1:A:119:ILE:HD12	1.84	0.58
1:B:158:GLU:OE2	1:B:291:HIS:O	2.20	0.58
1:C:199:THR:N	2:C:403:HOH:O	2.34	0.58
1:D:201:ALA:HA	1:D:204:HIS:HB3	1.86	0.58
1:A:268:VAL:O	1:A:272:ASN:ND2	2.37	0.57
1:B:196:GLU:HB2	1:B:203:TRP:CD2	2.39	0.57
1:D:201:ALA:O	1:D:204:HIS:HB3	2.04	0.57
1:D:210:LYS:HE3	1:D:213:GLU:HG3	1.85	0.57
1:C:209:ASN:OD1	1:C:211:MET:CB	2.52	0.57
1:C:80:ARG:HD2	1:C:122:TRP:CD2	2.39	0.57
1:D:266:SER:O	1:D:269:LYS:HB3	2.05	0.57
1:A:180:ARG:HB2	1:A:185:ILE:HD13	1.87	0.57
1:B:173:GLY:HA3	1:B:229:ILE:HG22	1.85	0.57
1:B:268:VAL:O	1:B:272:ASN:ND2	2.36	0.57
1:D:194:LEU:HA	1:D:195:ARG:HB2	1.86	0.56
1:D:224:LEU:HD23	1:D:250:ILE:HG12	1.87	0.56
1:D:35:VAL:HG23	1:D:91:SER:HB2	1.86	0.56
1:B:195:ARG:HH11	1:B:195:ARG:HG2	1.70	0.56
1:D:82:ARG:HG2	1:D:83:THR:N	2.20	0.56
1:A:296:ILE:HD12	1:A:297:LEU:N	2.20	0.56
1:A:268:VAL:HG12	1:A:272:ASN:ND2	2.19	0.56
1:B:82:ARG:NH1	1:B:337:LYS:HD2	2.21	0.56
1:C:85:ARG:NH1	1:C:334:GLU:OE2	2.38	0.56
1:D:156:GLN:OE1	1:D:156:GLN:N	2.38	0.56
1:A:108:GLN:CB	1:A:111:ASP:CG	2.62	0.56
1:B:134:ILE:H	1:B:134:ILE:HD12	1.70	0.56
1:A:29:VAL:O	1:A:32:TRP:N	2.33	0.56
1:D:4:ARG:HH22	1:D:121:LYS:CB	2.17	0.56
1:A:62:ARG:HH22	1:A:103:GLU:CD	2.06	0.56
1:B:292:THR:O	1:B:296:ILE:HG12	2.05	0.56
1:D:250:ILE:HB	1:D:259:LEU:HD22	1.88	0.56
1:D:194:LEU:C	1:D:203:TRP:HE1	2.09	0.55
1:C:209:ASN:OD1	1:C:211:MET:CG	2.53	0.55
1:B:320:ASP:O	1:B:322:LYS:HG3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:4:ARG:HH21	1:D:6:LEU:HD11	1.72	0.55
1:A:143:SER:O	1:A:225:GLN:NE2	2.40	0.55
1:A:139:PRO:HB2	1:A:142:ILE:HG13	1.88	0.55
1:B:240:LEU:HG	1:B:241:LYS:N	2.22	0.55
1:D:231:GLU:HA	1:D:236:ASN:ND2	2.22	0.55
1:A:147:GLN:O	1:A:326:LYS:HE2	2.07	0.54
1:A:161:LYS:CE	1:A:290:ASN:HA	2.36	0.54
1:C:164:LYS:NZ	1:C:288:GLU:OE1	2.41	0.54
1:D:231:GLU:HA	1:D:236:ASN:HD22	1.71	0.54
1:A:66:LEU:HD21	1:A:94:LEU:HD13	1.90	0.54
1:C:56:ILE:HG12	1:C:119:ILE:HD12	1.89	0.54
1:D:194:LEU:HB3	1:D:195:ARG:HB3	1.91	0.53
1:B:127:SER:HB2	1:B:130:LEU:HD12	1.90	0.53
1:B:337:LYS:HD3	2:B:409:HOH:O	2.08	0.53
1:B:85:ARG:O	1:B:86:LEU:HD23	2.09	0.53
1:B:80:ARG:HH12	1:B:338:ASP:HA	1.73	0.53
1:B:160:ASN:HB3	1:B:289:LEU:HD23	1.90	0.53
1:D:187:VAL:HB	1:D:203:TRP:NE1	2.24	0.53
1:D:82:ARG:NH2	1:D:337:LYS:CA	2.71	0.53
1:B:128:THR:O	1:B:131:SER:OG	2.15	0.53
1:A:133:GLU:HA	1:A:195:ARG:CD	2.39	0.53
1:A:196:GLU:HB3	1:A:203:TRP:CE2	2.43	0.53
1:C:155:PRO:O	1:C:158:GLU:HB2	2.09	0.53
1:C:0:SER:O	1:C:1:MET:O	2.27	0.52
1:A:302:GLY:O	1:A:311:ARG:HG2	2.10	0.52
1:B:157:ILE:C	1:B:158:GLU:HG2	2.19	0.52
1:D:96:LEU:HD22	1:D:100:ARG:HH21	1.73	0.52
1:A:30:ASP:HB3	1:A:154:ILE:HG21	1.92	0.52
1:B:59:ASP:HB3	1:B:63:TYR:HD2	1.74	0.52
1:D:338:ASP:N	1:D:338:ASP:OD1	2.43	0.52
1:B:63:TYR:OH	2:B:407:HOH:O	2.18	0.52
1:D:4:ARG:HH22	1:D:121:LYS:CG	2.23	0.52
1:B:287:LEU:HD13	1:B:288:GLU:C	2.30	0.51
1:D:204:HIS:CE1	1:D:208:ARG:HH21	2.28	0.51
1:B:161:LYS:HG3	1:B:290:ASN:HA	1.91	0.51
1:B:35:VAL:HG23	1:B:91:SER:HB2	1.92	0.51
1:D:9:ILE:HD11	1:D:104:LEU:HD13	1.91	0.51
1:D:37:LYS:O	1:D:40:GLU:HB2	2.10	0.51
1:D:130:LEU:HD21	1:D:201:ALA:HB3	1.92	0.51
1:B:237:LEU:HD22	1:B:332:TYR:HE2	1.76	0.51
1:D:65:MET:H	1:D:65:MET:HE3	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:ARG:NH1	1:B:103:GLU:OE2	2.44	0.51
1:D:82:ARG:NH2	1:D:337:LYS:CG	2.70	0.51
1:A:161:LYS:HD3	1:A:289:LEU:O	2.10	0.50
1:C:209:ASN:OD1	1:C:211:MET:HG2	2.11	0.50
1:B:207:ARG:HE	1:B:212:ILE:HD13	1.75	0.50
1:D:42:LYS:HD2	1:D:43:GLN:H	1.76	0.50
1:A:167:VAL:HG11	1:A:287:LEU:HD12	1.94	0.50
1:A:196:GLU:HB3	1:A:203:TRP:CD2	2.46	0.50
1:A:208:ARG:HG3	1:A:208:ARG:HH11	1.76	0.50
1:A:260:THR:HB	1:A:319:LEU:O	2.11	0.50
1:C:200:ASN:ND2	1:C:202:GLN:OE1	2.45	0.50
1:D:206:ALA:HA	1:D:211:MET:HE3	1.93	0.50
1:A:130:LEU:O	1:A:134:ILE:HG12	2.12	0.50
1:D:210:LYS:NZ	1:D:213:GLU:H	2.10	0.49
1:B:1:MET:N	2:B:402:HOH:O	2.04	0.49
1:C:336:HIS:O	1:C:339:ARG:HB2	2.12	0.49
1:A:307:LYS:HG2	1:A:307:LYS:O	2.11	0.49
1:C:170:LYS:HG2	1:C:312:GLU:OE2	2.12	0.49
1:A:303:PRO:O	1:A:311:ARG:NE	2.40	0.49
1:A:182:ASP:CG	1:C:338:ASP:OD2	2.50	0.49
1:C:241:LYS:HD3	1:C:241:LYS:H	1.76	0.49
1:D:296:ILE:HA	1:D:299:MET:HE3	1.94	0.49
1:C:68:LYS:HE3	2:C:410:HOH:O	2.12	0.49
1:D:4:ARG:CZ	1:D:4:ARG:HB3	2.43	0.49
1:A:3:VAL:HG12	1:A:4:ARG:N	2.26	0.49
1:C:116:ILE:HD12	1:C:116:ILE:H	1.78	0.49
1:D:326:LYS:HD2	1:D:328:ILE:HD11	1.94	0.48
1:B:237:LEU:HG	1:B:309:VAL:HG11	1.95	0.48
1:D:134:ILE:HG23	1:D:136:GLY:N	2.28	0.48
1:D:82:ARG:HH21	1:D:337:LYS:CA	2.26	0.48
1:C:318:ARG:NH2	1:C:320:ASP:OD2	2.46	0.48
1:C:339:ARG:HH11	1:C:339:ARG:CG	2.25	0.48
1:D:4:ARG:NH2	1:D:121:LYS:CG	2.76	0.48
1:A:128:THR:HG22	1:A:129:ASN:N	2.28	0.48
1:D:82:ARG:NH2	1:D:337:LYS:HB3	2.25	0.48
1:A:270:GLN:N	2:A:401:HOH:O	2.42	0.48
1:B:115:GLU:OE2	2:B:408:HOH:O	2.20	0.48
1:C:85:ARG:HD2	1:C:334:GLU:OE1	2.14	0.48
1:A:262:ILE:O	1:A:266:SER:N	2.38	0.48
1:A:62:ARG:NH2	1:A:101:PHE:HD1	2.11	0.48
1:D:178:VAL:HA	1:D:186:GLY:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:54:SER:OG	1:D:81:LEU:HD12	2.13	0.48
1:D:62:ARG:HD3	1:D:101:PHE:CE2	2.49	0.48
1:D:190:ARG:NH1	1:D:191:ASN:CG	2.66	0.48
1:D:197:THR:HG22	1:D:199:THR:H	1.79	0.48
1:B:226:GLY:HA3	1:B:246:TYR:O	2.14	0.47
1:D:255:LYS:HB2	1:D:257:GLN:HG2	1.96	0.47
1:D:261:PRO:HG2	1:D:319:LEU:HA	1.96	0.47
1:A:108:GLN:CD	1:A:111:ASP:OD1	2.52	0.47
1:B:37:LYS:O	1:B:40:GLU:HB2	2.14	0.47
1:D:229:ILE:O	1:D:243:GLN:HA	2.15	0.47
1:D:223:ALA:O	1:D:250:ILE:HA	2.14	0.47
1:C:85:ARG:HH11	1:C:334:GLU:CD	2.17	0.47
1:B:237:LEU:O	1:B:237:LEU:HG	2.14	0.47
1:D:121:LYS:HD3	1:D:123:GLU:HG3	1.95	0.47
1:A:259:LEU:HB3	1:A:263:GLU:HB2	1.96	0.47
1:A:233:ILE:CA	1:A:234:GLN:OE1	2.60	0.47
1:A:160:ASN:O	1:A:289:LEU:HD23	2.15	0.47
1:D:190:ARG:HH11	1:D:190:ARG:HB3	1.80	0.47
1:B:320:ASP:CG	1:B:322:LYS:HB2	2.35	0.47
1:C:58:GLU:OE1	1:C:77:ARG:HD3	2.14	0.47
1:A:150:VAL:HG11	1:A:296:ILE:CD1	2.45	0.47
1:A:161:LYS:CD	1:A:289:LEU:O	2.63	0.47
1:C:133:GLU:N	1:C:133:GLU:OE1	2.48	0.47
1:C:318:ARG:HH21	1:C:322:LYS:HB2	1.79	0.47
1:D:209:ASN:OD1	1:D:211:MET:HG3	2.15	0.46
1:A:208:ARG:CG	1:A:208:ARG:HH11	2.28	0.46
1:A:268:VAL:CG1	1:A:272:ASN:HD21	2.27	0.46
1:A:83:THR:HG22	1:A:92:GLN:HG2	1.96	0.46
1:A:150:VAL:HG21	1:A:296:ILE:CD1	2.46	0.46
1:C:209:ASN:OD1	1:C:211:MET:HB3	2.15	0.46
1:B:185:ILE:HG21	1:B:212:ILE:HG23	1.97	0.46
1:D:135:LEU:HD11	1:D:195:ARG:H	1.81	0.46
1:D:164:LYS:HA	1:D:287:LEU:O	2.16	0.46
1:A:108:GLN:HB2	1:A:111:ASP:OD1	2.11	0.46
1:A:176:MET:HA	1:A:188:CYS:O	2.16	0.46
1:B:319:LEU:HA	1:B:319:LEU:HD23	1.63	0.45
1:A:304:SER:HB2	1:A:311:ARG:HA	1.99	0.45
1:D:179:TYR:O	1:D:185:ILE:HA	2.17	0.45
1:A:247:LEU:HG	1:A:248:PHE:N	2.31	0.45
1:A:83:THR:HG21	1:A:333:LEU:HD13	1.99	0.45
1:B:144:LYS:HE2	1:B:190:ARG:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:PRO:O	1:A:159:GLU:HG3	2.17	0.45
1:C:154:ILE:HB	1:C:155:PRO:HD3	1.99	0.45
1:D:233:ILE:N	1:D:233:ILE:HD12	2.32	0.45
1:B:261:PRO:HG2	1:B:319:LEU:HD23	1.99	0.44
1:D:230:GLY:HA3	1:D:243:GLN:HG3	1.99	0.44
1:D:259:LEU:HA	1:D:259:LEU:HD12	1.76	0.44
1:D:303:PRO:O	1:D:311:ARG:HD3	2.18	0.44
1:A:317:LYS:HD3	1:A:324:SER:HB3	1.98	0.44
1:A:182:ASP:HB3	1:C:338:ASP:OD2	2.17	0.44
1:A:128:THR:HG21	1:B:134:ILE:HB	2.00	0.44
1:D:300:ALA:HA	1:D:314:LEU:HG	1.99	0.44
1:D:82:ARG:CD	1:D:82:ARG:C	2.85	0.44
1:A:328:ILE:H	1:A:328:ILE:HD12	1.82	0.44
1:C:157:ILE:HD13	1:C:157:ILE:HA	1.79	0.44
1:D:190:ARG:HH12	1:D:191:ASN:ND2	2.14	0.44
1:A:250:ILE:HD13	1:A:267:LEU:HD23	2.00	0.44
1:C:29:VAL:O	1:C:32:TRP:N	2.43	0.44
1:A:104:LEU:HA	1:A:107:ASN:OD1	2.18	0.44
1:A:317:LYS:HB3	1:A:324:SER:CB	2.35	0.44
1:C:312:GLU:HG3	1:C:332:TYR:CE1	2.52	0.44
1:A:6:LEU:HD11	1:A:121:LYS:HB2	2.00	0.43
1:A:214:GLY:O	1:A:218:LEU:HG	2.18	0.43
1:D:4:ARG:NH2	1:D:121:LYS:HG3	2.33	0.43
1:D:42:LYS:HA	1:D:42:LYS:HD2	1.71	0.43
1:A:180:ARG:CB	1:A:185:ILE:HD13	2.48	0.43
1:A:246:TYR:CE2	1:A:305:LEU:HD11	2.53	0.43
1:D:68:LYS:HD2	1:D:68:LYS:HA	1.88	0.43
1:A:310:LYS:CD	1:A:335:LYS:HD2	2.48	0.43
1:C:169:VAL:HA	1:C:314:LEU:HD23	2.00	0.43
1:D:174:SER:O	1:D:228:ILE:HG22	2.18	0.43
1:B:231:GLU:O	1:B:243:GLN:NE2	2.48	0.43
1:C:198:ALA:N	2:C:403:HOH:O	2.50	0.43
1:C:34:CYS:SG	2:C:425:HOH:O	2.62	0.43
1:D:264:ARG:O	1:D:268:VAL:HG12	2.18	0.43
1:D:59:ASP:HB3	1:D:63:TYR:HD2	1.82	0.43
1:C:268:VAL:HG21	1:C:280:HIS:HD2	1.83	0.43
1:D:82:ARG:HD3	1:D:82:ARG:C	2.38	0.43
1:B:20:GLY:O	1:B:87:ARG:NH2	2.52	0.43
1:B:137:GLU:N	1:B:137:GLU:CD	2.72	0.43
1:B:82:ARG:HH11	1:B:337:LYS:HD2	1.82	0.43
1:B:59:ASP:HB3	1:B:63:TYR:CD2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:207:ARG:HE	1:D:210:LYS:HE2	1.83	0.43
1:D:240:LEU:HG	1:D:241:LYS:N	2.34	0.43
1:A:246:TYR:CZ	1:A:305:LEU:HD11	2.54	0.42
1:B:197:THR:HB	1:B:199:THR:HG22	2.00	0.42
1:B:97:PRO:HB2	1:B:99:ASP:OD1	2.18	0.42
1:A:113:VAL:HG12	1:A:116:ILE:HD12	2.01	0.42
1:A:249:ASP:HB3	1:A:258:TYR:CE2	2.54	0.42
1:A:269:LYS:O	1:A:273:ASP:N	2.35	0.42
1:A:55:PHE:CZ	1:A:73:TYR:HB3	2.54	0.42
1:B:217:PHE:CZ	1:D:72:ASP:HB2	2.53	0.42
1:B:46:ARG:HG3	1:B:98:MET:HG3	2.01	0.42
1:A:180:ARG:HB2	1:A:185:ILE:CD1	2.50	0.42
1:A:150:VAL:HG21	1:A:296:ILE:HD11	2.01	0.42
1:D:184:HIS:NE2	1:D:193:GLU:OE1	2.44	0.42
1:D:201:ALA:HA	1:D:204:HIS:CB	2.49	0.42
1:D:171:LEU:HD12	1:D:281:VAL:HG12	2.02	0.42
1:B:96:LEU:HA	1:B:96:LEU:HD23	1.81	0.42
1:D:4:ARG:HH22	1:D:121:LYS:HG3	1.84	0.42
1:C:218:LEU:HD23	1:C:218:LEU:HA	1.78	0.42
1:D:187:VAL:HB	1:D:203:TRP:CE2	2.55	0.42
1:D:43:GLN:O	2:D:402:HOH:O	2.21	0.42
1:B:137:GLU:CD	1:B:137:GLU:H	2.23	0.42
1:A:247:LEU:HD21	1:A:250:ILE:HG13	2.01	0.42
1:C:223:ALA:HB3	1:C:251:TYR:HB3	2.01	0.42
1:C:250:ILE:HD13	1:C:267:LEU:HD22	2.02	0.42
1:D:265:GLN:NE2	2:D:403:HOH:O	2.37	0.42
1:D:293:ALA:HA	1:D:296:ILE:HD12	2.02	0.42
1:C:336:HIS:HB3	1:C:339:ARG:HG3	2.02	0.41
1:B:147:GLN:HB2	1:B:326:LYS:HG2	2.02	0.41
1:A:19:ALA:C	1:A:21:ALA:H	2.24	0.41
1:A:142:ILE:HA	1:A:251:TYR:CE2	2.56	0.41
1:B:85:ARG:NH1	1:B:334:GLU:OE1	2.53	0.41
1:C:37:LYS:O	1:C:40:GLU:HB2	2.19	0.41
1:D:259:LEU:HD12	1:D:263:GLU:OE1	2.20	0.41
1:A:180:ARG:HD3	1:A:220:ARG:O	2.20	0.41
1:B:55:PHE:CZ	1:B:73:TYR:HB3	2.55	0.41
1:C:287:LEU:HD12	1:C:288:GLU:N	2.35	0.41
1:A:217:PHE:CZ	1:C:72:ASP:HB2	2.55	0.41
1:A:161:LYS:HZ1	1:A:291:HIS:N	2.18	0.41
1:B:58:GLU:OE1	1:B:77:ARG:HD3	2.20	0.41
1:C:65:MET:N	1:C:65:MET:SD	2.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:139:PRO:HB3	1:D:141:PHE:CE1	2.56	0.41
1:D:230:GLY:HA3	1:D:233:ILE:HD13	2.02	0.41
1:C:172:ASP:HB2	1:C:237:LEU:HD21	2.03	0.41
1:D:177:THR:OG1	1:D:188:CYS:HB2	2.20	0.41
1:A:247:LEU:HD23	1:A:280:HIS:HD2	1.85	0.41
1:B:80:ARG:HD2	1:B:122:TRP:CD2	2.55	0.41
1:C:85:ARG:NH1	1:C:334:GLU:CD	2.74	0.41
1:D:226:GLY:HA3	1:D:246:TYR:O	2.21	0.41
1:D:334:GLU:O	1:D:337:LYS:HG3	2.21	0.41
1:C:264:ARG:O	1:C:268:VAL:HG23	2.21	0.41
1:D:139:PRO:HB3	1:D:141:PHE:CZ	2.55	0.41
1:A:312:GLU:HG3	1:A:332:TYR:CE1	2.55	0.41
1:B:69:GLN:HG3	1:B:79:THR:CG2	2.51	0.41
1:A:172:ASP:OD2	1:A:237:LEU:HD23	2.21	0.40
1:C:35:VAL:HG23	1:C:91:SER:HB2	2.02	0.40
1:C:80:ARG:HH22	1:C:339:ARG:N	2.19	0.40
1:A:108:GLN:OE1	1:A:111:ASP:OD1	2.39	0.40
1:A:128:THR:HG22	1:A:129:ASN:H	1.86	0.40
1:B:37:LYS:HE3	1:B:37:LYS:HB3	1.52	0.40
1:D:130:LEU:HA	1:D:130:LEU:HD23	1.89	0.40
1:D:46:ARG:HG2	1:D:46:ARG:H	1.68	0.40
1:A:46:ARG:O	1:A:98:MET:HG2	2.22	0.40
1:B:157:ILE:HG22	1:B:158:GLU:HG2	2.02	0.40
1:C:226:GLY:HA2	1:C:248:PHE:CE2	2.57	0.40
1:D:297:LEU:HD23	1:D:297:LEU:HA	1.91	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)
1:A:72:ASP:OD2	1:D:270:GLN:NE2[4_555]	2.17	0.03
1:A:77:ARG:NH1	1:D:274:ASN:OD1[4_555]	2.19	0.01

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	338/340 (99%)	327 (97%)	11 (3%)	0	100	100
1	B	338/340 (99%)	327 (97%)	10 (3%)	1 (0%)	41	61
1	C	338/340 (99%)	323 (96%)	13 (4%)	2 (1%)	25	43
1	D	338/340 (99%)	323 (96%)	13 (4%)	2 (1%)	25	43
All	All	1352/1360 (99%)	1300 (96%)	47 (4%)	5 (0%)	34	54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1	MET
1	B	30	ASP
1	D	195	ARG
1	C	30	ASP
1	D	30	ASP

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	293/293 (100%)	282 (96%)	11 (4%)	33	58
1	B	293/293 (100%)	285 (97%)	8 (3%)	44	71
1	C	293/293 (100%)	284 (97%)	9 (3%)	40	67
1	D	293/293 (100%)	282 (96%)	11 (4%)	33	58
All	All	1172/1172 (100%)	1133 (97%)	39 (3%)	38	64

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	LYS
1	A	46	ARG
1	A	115	GLU

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Mol	Chain	Res	Type
1	A	127	SER
1	A	161	LYS
1	A	175	SER
1	A	204	HIS
1	A	209	ASN
1	A	248	PHE
1	A	322	LYS
1	A	323	PHE
1	B	22	GLU
1	B	34	CYS
1	B	46	ARG
1	B	164	LYS
1	B	195	ARG
1	B	207	ARG
1	B	208	ARG
1	B	266	SER
1	C	61	ASP
1	C	82	ARG
1	C	135	LEU
1	C	174	SER
1	C	209	ASN
1	C	241	LYS
1	C	248	PHE
1	C	266	SER
1	C	310	LYS
1	D	34	CYS
1	D	40	GLU
1	D	82	ARG
1	D	174	SER
1	D	181	LYS
1	D	183	ASP
1	D	204	HIS
1	D	239	LYS
1	D	291	HIS
1	D	322	LYS
1	D	338	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	272	ASN

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Mol	Chain	Res	Type
1	B	108	GLN
1	D	202	GLN
1	D	204	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	340/340 (100%)	0.55	22 (6%) 18 19	50, 91, 124, 173	0
1	B	340/340 (100%)	0.16	11 (3%) 47 51	35, 70, 109, 149	0
1	C	340/340 (100%)	0.15	10 (2%) 51 55	34, 61, 103, 134	0
1	D	340/340 (100%)	0.83	38 (11%) 5 4	55, 99, 139, 161	0
All	All	1360/1360 (100%)	0.42	81 (5%) 21 22	34, 80, 126, 173	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	2	GLY	16.6
1	A	0	SER	10.3
1	B	2	GLY	10.1
1	C	0	SER	7.3
1	A	2	GLY	6.7
1	A	1	MET	6.5
1	D	240	LEU	5.3
1	A	290	ASN	5.3
1	D	237	LEU	5.1
1	D	130	LEU	5.1
1	D	1	MET	5.0
1	B	1	MET	4.9
1	D	132	GLY	4.7
1	D	135	LEU	4.7
1	D	3	VAL	4.6
1	D	289	LEU	4.4
1	A	104	LEU	4.2
1	D	0	SER	4.2
1	A	3	VAL	4.2
1	B	0	SER	4.2
1	D	198	ALA	4.0

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Mol	Chain	Res	Type	RSRZ
1	D	323	PHE	3.9
1	C	339	ARG	3.9
1	D	141	PHE	3.9
1	B	339	ARG	3.9
1	C	4	ARG	3.9
1	D	239	LYS	3.8
1	A	289	LEU	3.7
1	C	34	CYS	3.7
1	B	181	LYS	3.6
1	D	129	ASN	3.6
1	D	137	GLU	3.4
1	A	307	LYS	3.4
1	A	135	LEU	3.4
1	D	19	ALA	3.2
1	A	308	ASN	3.2
1	B	3	VAL	3.2
1	D	128	THR	3.2
1	C	235	GLY	3.1
1	D	131	SER	3.1
1	D	190	ARG	3.1
1	D	241	LYS	3.0
1	D	136	GLY	3.0
1	D	288	GLU	3.0
1	A	328	ILE	2.9
1	D	17	PRO	2.9
1	C	240	LEU	2.8
1	B	338	ASP	2.8
1	A	127	SER	2.8
1	A	192	TRP	2.8
1	B	192	TRP	2.8
1	D	254	ASP	2.8
1	A	126	ILE	2.7
1	C	129	ASN	2.7
1	D	305	LEU	2.7
1	A	330	ASN	2.7
1	D	219	ASN	2.6
1	D	185	ILE	2.6
1	C	1	MET	2.6
1	D	291	HIS	2.5
1	D	308	ASN	2.5
1	A	284	LEU	2.5
1	A	157	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	14	GLU	2.4
1	D	307	LYS	2.3
1	A	240	LEU	2.3
1	B	135	LEU	2.3
1	B	158	GLU	2.3
1	B	240	LEU	2.3
1	A	108	GLN	2.2
1	C	237	LEU	2.2
1	D	278	VAL	2.2
1	D	228	ILE	2.2
1	D	181	LYS	2.2
1	A	241	LYS	2.2
1	D	243	GLN	2.2
1	D	21	ALA	2.2
1	A	162	GLY	2.2
1	A	195	ARG	2.2
1	C	127	SER	2.2
1	D	159	GLU	2.1

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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.