



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:24 PM UTC

PDB ID : 3UXI / pdb_00003uxi
Title : Crystal structure of L-rhamnose isomerase W38A mutant from *Bacillus halodurans*
Authors : Doan, T.T.N.; Prabhu, P.; Kim, J.K.; Jeya, M.; Kang, L.W.; Lee, J.K.
Deposited on : 2011-12-05
Resolution : 2.73 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

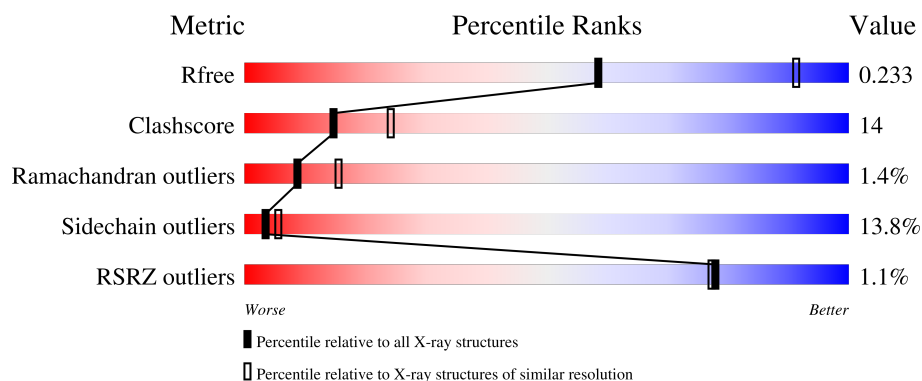
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

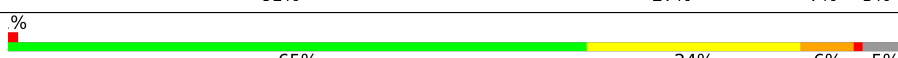
The reported resolution of this entry is 2.73 Å.

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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	424	
1	B	424	
1	C	424	
1	D	424	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13410 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 L-Rhamnose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	0	0
			3278	2098	564	607	9			
1	B	402	Total	C	N	O	S	0	0	0
			3278	2098	564	607	9			
1	C	404	Total	C	N	O	S	0	0	0
			3295	2109	567	609	10			
1	D	404	Total	C	N	O	S	0	0	0
			3295	2109	567	609	10			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP Q9KCL9
A	-4	HIS	-	expression tag	UNP Q9KCL9
A	-3	HIS	-	expression tag	UNP Q9KCL9
A	-2	HIS	-	expression tag	UNP Q9KCL9
A	-1	HIS	-	expression tag	UNP Q9KCL9
A	0	HIS	-	expression tag	UNP Q9KCL9
A	38	ALA	TRP	engineered mutation	UNP Q9KCL9
B	-5	HIS	-	expression tag	UNP Q9KCL9
B	-4	HIS	-	expression tag	UNP Q9KCL9
B	-3	HIS	-	expression tag	UNP Q9KCL9
B	-2	HIS	-	expression tag	UNP Q9KCL9
B	-1	HIS	-	expression tag	UNP Q9KCL9
B	0	HIS	-	expression tag	UNP Q9KCL9
B	38	ALA	TRP	engineered mutation	UNP Q9KCL9
C	-5	HIS	-	expression tag	UNP Q9KCL9
C	-4	HIS	-	expression tag	UNP Q9KCL9
C	-3	HIS	-	expression tag	UNP Q9KCL9
C	-2	HIS	-	expression tag	UNP Q9KCL9
C	-1	HIS	-	expression tag	UNP Q9KCL9
C	0	HIS	-	expression tag	UNP Q9KCL9
C	38	ALA	TRP	engineered mutation	UNP Q9KCL9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-5	HIS	-	expression tag	UNP Q9KCL9
D	-4	HIS	-	expression tag	UNP Q9KCL9
D	-3	HIS	-	expression tag	UNP Q9KCL9
D	-2	HIS	-	expression tag	UNP Q9KCL9
D	-1	HIS	-	expression tag	UNP Q9KCL9
D	0	HIS	-	expression tag	UNP Q9KCL9
D	38	ALA	TRP	engineered mutation	UNP Q9KCL9

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mn 1 1	0	0

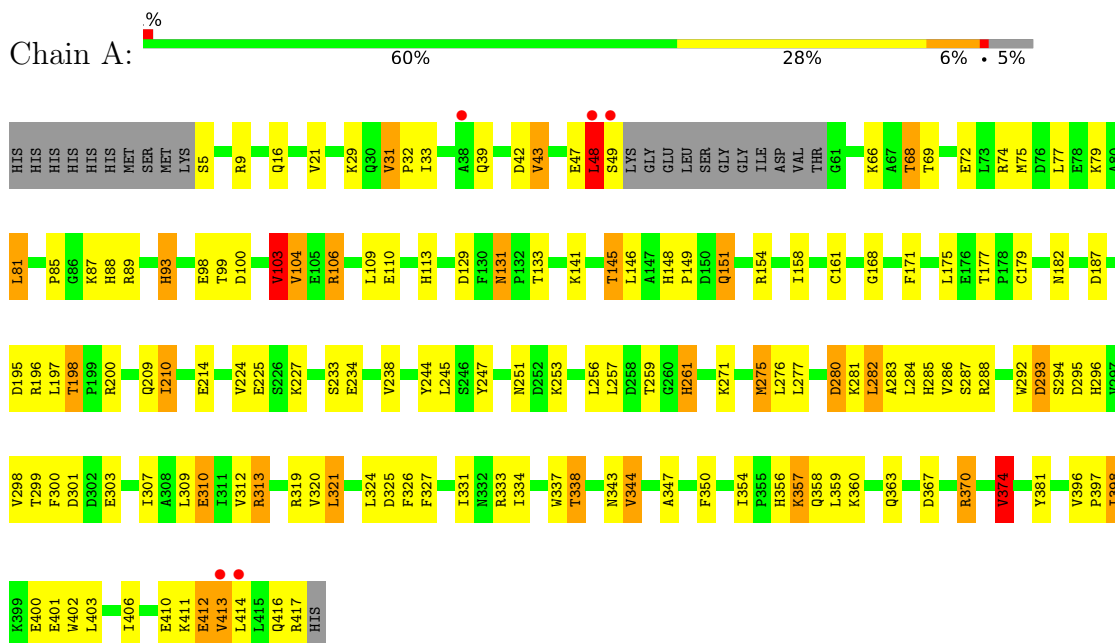
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	76	Total O 76 76	0	0
3	B	62	Total O 62 62	0	0
3	C	56	Total O 56 56	0	0
3	D	69	Total O 69 69	0	0

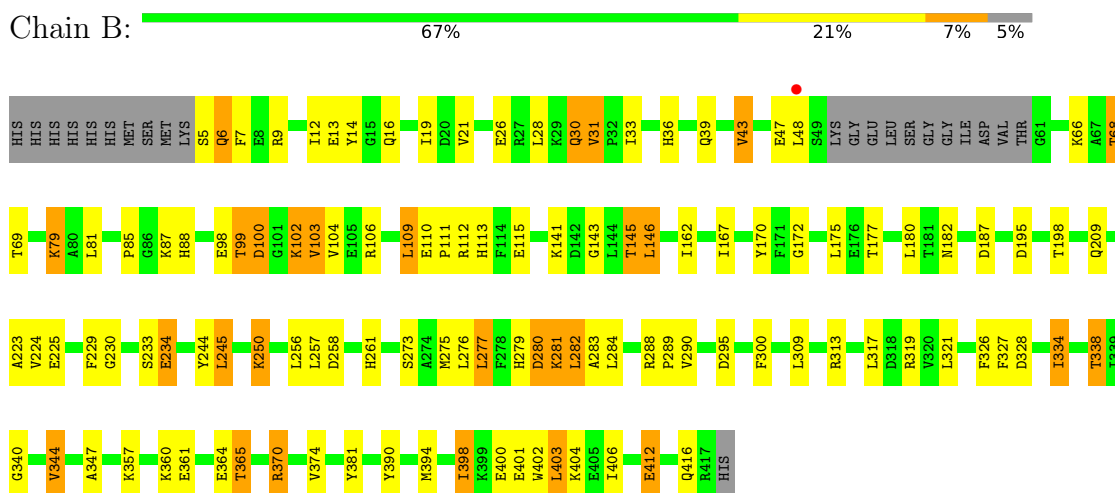
3 Residue-property plots

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.

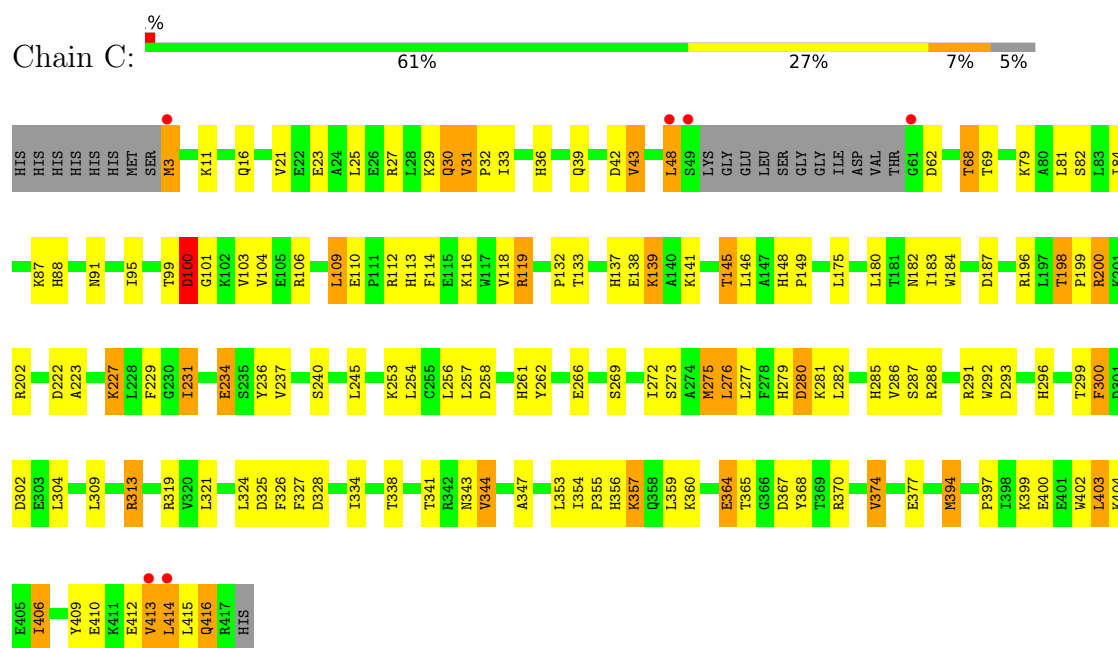
• Molecule 1: L-Rhamnose isomerase



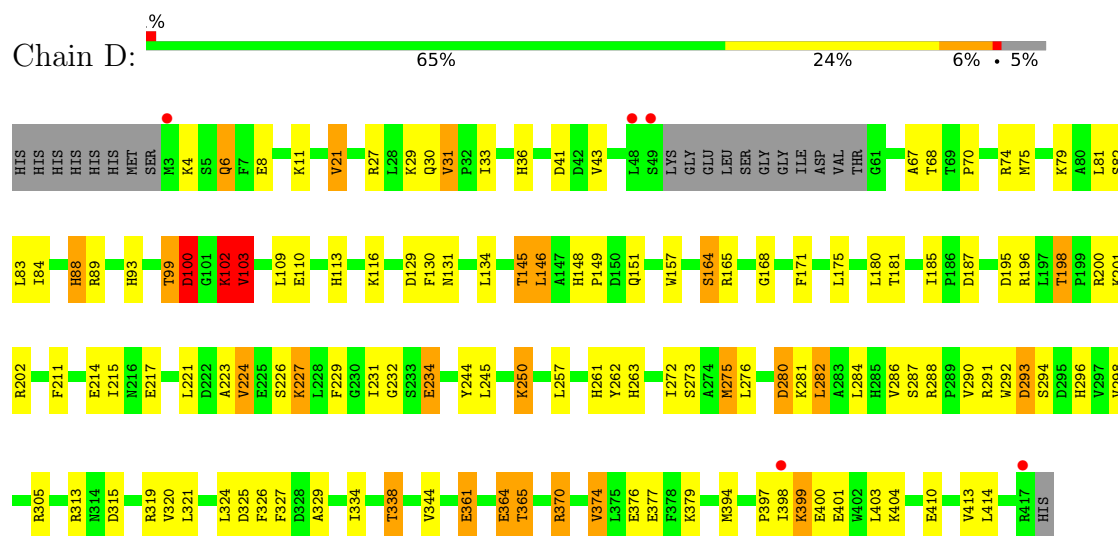
• Molecule 1: L-Rhamnose isomerase



• Molecule 1: L-Rhamnose isomerase



- Molecule 1: L-Rhamnose isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.38Å 164.79Å 92.84Å 90.00° 117.02° 90.00°	Depositor
Resolution (Å)	45.87 – 2.73 45.87 – 2.73	Depositor EDS
% Data completeness (in resolution range)	82.5 (45.87-2.73) 82.5 (45.87-2.73)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.186 , 0.267 (Not available) , 0.233	Depositor DCC
R_{free} test set	2480 reflections (4.19%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 27.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13410	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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: MN

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	1.01	6/3357 (0.2%)	1.13	9/4545 (0.2%)
1	B	0.86	0/3357	1.12	10/4545 (0.2%)
1	C	0.94	2/3374 (0.1%)	1.07	2/4566 (0.0%)
1	D	0.88	1/3374 (0.0%)	1.13	9/4566 (0.2%)
All	All	0.92	9/13462 (0.1%)	1.11	30/18222 (0.2%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	261	HIS	CG-ND1	-7.79	1.29	1.38
1	A	261	HIS	CD2-NE2	-6.96	1.30	1.37
1	C	198	THR	C-O	-6.45	1.18	1.24
1	D	103	VAL	CA-CB	6.37	1.63	1.54
1	C	200	ARG	C-O	-5.69	1.17	1.24
1	A	48	LEU	CA-C	5.27	1.59	1.52
1	A	259	THR	C-O	-5.20	1.17	1.24
1	A	257	LEU	C-O	-5.11	1.17	1.24
1	A	103	VAL	CA-CB	5.01	1.61	1.54

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	398	ILE	N-CA-C	11.55	123.82	111.58
1	A	99	THR	N-CA-C	-11.34	96.46	110.41
1	B	30	GLN	CA-C-N	-11.20	114.53	122.59
1	B	30	GLN	C-N-CA	-11.20	114.53	122.59
1	A	374	VAL	N-CA-C	8.76	118.83	110.42
1	B	284	LEU	N-CA-C	8.06	121.72	109.23
1	B	102	LYS	N-CA-C	7.86	119.63	111.14
1	A	31	VAL	N-CA-C	7.54	115.11	108.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ARG	N-CA-C	7.28	121.06	111.75
1	B	381	TYR	CA-C-N	6.75	128.28	119.84
1	B	381	TYR	C-N-CA	6.75	128.28	119.84
1	D	146	LEU	N-CA-C	-6.62	105.17	113.18
1	D	313	ARG	CB-CA-C	-6.58	99.49	110.68
1	D	374	VAL	N-CA-C	6.28	117.84	111.00
1	D	398	ILE	N-CA-C	6.23	117.72	110.62
1	D	102	LYS	N-CA-C	5.71	120.38	113.41
1	B	99	THR	N-CA-C	-5.59	98.90	110.80
1	B	79	LYS	N-CA-C	-5.58	105.10	111.07
1	A	177	THR	CA-C-N	-5.45	113.95	119.83
1	A	177	THR	C-N-CA	-5.45	113.95	119.83
1	C	262	TYR	N-CA-C	5.44	117.53	109.69
1	D	88	HIS	N-CA-C	5.43	118.69	108.65
1	D	262	TYR	N-CA-C	5.41	118.09	109.81
1	B	146	LEU	N-CA-C	-5.39	106.66	113.18
1	B	398	ILE	N-CA-C	5.31	117.21	111.58
1	A	261	HIS	N-CA-C	5.28	118.74	112.72
1	D	31	VAL	CB-CA-C	5.26	115.83	110.94
1	D	272	ILE	N-CA-C	5.12	115.56	110.23
1	C	99	THR	N-CA-C	-5.09	103.32	110.35
1	A	48	LEU	N-CA-C	5.04	121.54	110.80

There are no chirality outliers.

There are no planarity outliers.

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	3278	0	3231	105	0
1	B	3278	0	3231	79	0
1	C	3295	0	3253	106	0
1	D	3295	0	3253	99	0
2	A	1	0	0	0	0
3	A	76	0	0	3	0
3	B	62	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	56	0	0	2	0
3	D	69	0	0	2	0
All	All	13410	0	12968	364	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 14.

All (364) 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:195:ASP:OD2	1:D:198:THR:HG23	1.50	1.12
1:A:410:GLU:HA	1:A:414:LEU:HB2	1.35	1.06
1:B:195:ASP:OD2	1:B:198:THR:HG22	1.55	1.05
1:C:245:LEU:HD22	1:D:245:LEU:HD22	1.43	1.00
1:D:286:VAL:HG23	1:D:298:VAL:HG21	1.46	0.95
1:C:110:GLU:H	1:C:113:HIS:HD2	1.13	0.95
1:C:245:LEU:HD22	1:D:245:LEU:CD2	1.96	0.94
1:A:286:VAL:HG23	1:A:298:VAL:HG21	1.51	0.92
1:D:287:SER:HB3	1:D:325:ASP:O	1.68	0.92
1:B:370:ARG:HG3	1:B:370:ARG:HH11	1.32	0.92
1:A:5:SER:HB2	3:A:458:HOH:O	1.70	0.91
1:D:148:HIS:ND1	1:D:149:PRO:HD2	1.88	0.89
1:C:257:LEU:HD11	1:C:275:MET:HE1	1.55	0.88
1:A:110:GLU:H	1:A:113:HIS:HD2	1.19	0.88
1:C:145:THR:HG22	1:C:146:LEU:H	1.35	0.88
1:C:280:ASP:HA	1:C:319:ARG:NH1	1.89	0.87
1:D:195:ASP:OD2	1:D:198:THR:CG2	2.23	0.86
1:B:145:THR:CG2	1:B:146:LEU:H	1.89	0.85
1:A:261:HIS:CD2	1:A:293:ASP:OD2	2.28	0.85
1:C:280:ASP:HA	1:C:319:ARG:HH11	1.39	0.85
1:D:110:GLU:H	1:D:113:HIS:HD2	1.25	0.84
1:B:334:ILE:O	1:B:338:THR:HG23	1.77	0.84
1:D:370:ARG:HH11	1:D:370:ARG:HG3	1.44	0.82
1:A:85:PRO:O	1:A:88:HIS:HE1	1.60	0.82
1:A:334:ILE:O	1:A:338:THR:HG23	1.79	0.82
1:C:145:THR:CG2	1:C:146:LEU:H	1.92	0.82
1:D:334:ILE:O	1:D:338:THR:HG22	1.78	0.82
1:D:145:THR:HG22	1:D:146:LEU:N	1.96	0.81
1:B:195:ASP:CG	1:B:198:THR:HG22	2.05	0.81
1:B:145:THR:HG22	1:B:146:LEU:H	1.46	0.80
1:C:110:GLU:H	1:C:113:HIS:CD2	1.99	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:410:GLU:HA	1:C:414:LEU:HB2	1.61	0.80
1:B:110:GLU:H	1:B:113:HIS:HD2	1.27	0.79
1:D:134:LEU:HD23	1:D:145:THR:HG21	1.65	0.79
1:A:195:ASP:OD2	1:A:198:THR:CG2	2.31	0.79
1:B:145:THR:CG2	1:B:146:LEU:N	2.45	0.79
1:A:68:THR:HG22	1:A:69:THR:HG23	1.66	0.77
1:A:68:THR:HB	1:A:72:GLU:OE1	1.85	0.77
1:C:137:HIS:HE1	1:C:139:LYS:HG3	1.50	0.76
1:A:145:THR:HG22	1:A:146:LEU:H	1.49	0.76
1:D:257:LEU:HD11	1:D:275:MET:CE	2.14	0.76
1:D:370:ARG:HH11	1:D:370:ARG:CG	1.97	0.76
1:D:280:ASP:O	1:D:319:ARG:HD2	1.87	0.74
1:D:334:ILE:O	1:D:338:THR:CG2	2.34	0.74
1:C:145:THR:CG2	1:C:146:LEU:N	2.50	0.74
1:B:280:ASP:HA	1:B:319:ARG:NH1	2.02	0.74
1:A:110:GLU:H	1:A:113:HIS:CD2	2.05	0.73
1:A:47:GLU:O	1:A:49:SER:N	2.23	0.72
1:D:33:ILE:H	1:D:88:HIS:HD2	1.35	0.72
1:C:137:HIS:CE1	1:C:139:LYS:HG3	2.25	0.71
1:A:145:THR:HB	1:A:187:ASP:OD1	1.91	0.71
1:C:198:THR:HB	1:C:199:PRO:HD3	1.72	0.71
1:C:198:THR:HB	1:C:199:PRO:CD	2.21	0.71
1:A:287:SER:HB3	1:A:325:ASP:O	1.92	0.69
1:A:145:THR:HG22	1:A:146:LEU:N	2.03	0.69
1:A:295:ASP:HB3	1:A:327:PHE:H	1.56	0.69
1:D:130:PHE:CE2	1:D:164:SER:HB3	2.27	0.69
1:C:132:PRO:HD2	1:C:182:ASN:O	1.92	0.69
1:B:110:GLU:H	1:B:113:HIS:CD2	2.08	0.69
1:D:130:PHE:HE2	1:D:164:SER:HB3	1.58	0.68
1:A:33:ILE:H	1:A:88:HIS:CD2	2.12	0.68
1:B:145:THR:HG23	1:B:146:LEU:N	2.09	0.68
1:B:102:LYS:O	1:B:103:VAL:HG12	1.94	0.67
1:C:79:LYS:O	1:C:82:SER:HB2	1.94	0.67
1:A:299:THR:HA	1:A:343:ASN:HD22	1.58	0.66
1:C:196:ARG:HG2	1:C:236:TYR:CE1	2.31	0.66
1:B:280:ASP:O	1:B:319:ARG:HD2	1.97	0.65
1:A:148:HIS:ND1	1:A:149:PRO:HD2	2.10	0.65
1:B:195:ASP:OD2	1:B:198:THR:CG2	2.37	0.65
1:B:370:ARG:HH11	1:B:370:ARG:CG	2.08	0.65
1:A:234:GLU:OE2	1:D:292:TRP:NE1	2.26	0.65
1:D:148:HIS:CE1	1:D:149:PRO:HD2	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:ASP:HA	1:A:319:ARG:NH1	2.13	0.64
1:A:33:ILE:H	1:A:88:HIS:HD2	1.42	0.64
1:A:261:HIS:NE2	1:A:293:ASP:OD2	2.31	0.64
1:D:130:PHE:O	1:D:181:THR:HA	1.96	0.64
1:D:145:THR:HB	1:D:187:ASP:OD1	1.97	0.64
1:A:271:LYS:HB3	1:A:275:MET:HE1	1.77	0.64
1:B:280:ASP:HA	1:B:319:ARG:HH11	1.61	0.63
1:C:180:LEU:HD11	1:C:223:ALA:HB2	1.80	0.63
1:D:227:LYS:HA	1:D:261:HIS:HB3	1.80	0.63
1:A:85:PRO:O	1:A:88:HIS:CE1	2.49	0.63
1:D:324:LEU:HD13	1:D:326:PHE:CE2	2.33	0.62
1:B:279:HIS:O	1:B:280:ASP:CB	2.47	0.62
1:A:288:ARG:HB3	1:A:296:HIS:HB2	1.81	0.62
1:C:222:ASP:HB2	1:C:253:LYS:HD2	1.82	0.62
1:B:279:HIS:O	1:B:280:ASP:HB3	2.00	0.61
1:C:256:LEU:HD21	1:C:285:HIS:CG	2.36	0.61
1:C:231:ILE:O	1:C:231:ILE:HG13	2.00	0.61
1:C:356:HIS:O	1:C:360:LYS:HG3	2.01	0.61
1:D:324:LEU:HD13	1:D:326:PHE:HE2	1.65	0.61
1:A:292:TRP:O	1:A:294:SER:N	2.34	0.60
1:D:327:PHE:CZ	1:D:329:ALA:HB2	2.37	0.60
1:C:280:ASP:O	1:C:319:ARG:HD2	2.01	0.60
1:A:148:HIS:CG	1:A:149:PRO:HD2	2.37	0.59
1:C:39:GLN:HE22	1:C:328:ASP:H	1.50	0.59
1:A:195:ASP:OD2	1:A:198:THR:HG23	2.01	0.59
1:C:148:HIS:ND1	1:C:149:PRO:HD2	2.17	0.59
1:A:296:HIS:CE1	1:D:234:GLU:HG3	2.37	0.59
1:B:257:LEU:HD11	1:B:275:MET:CE	2.33	0.59
1:D:361:GLU:O	1:D:365:THR:HB	2.03	0.59
1:C:394:MET:HA	1:C:394:MET:HE2	1.85	0.58
1:A:48:LEU:HD13	1:A:48:LEU:N	2.19	0.58
1:B:6:GLN:HG2	1:B:7:PHE:N	2.18	0.58
1:A:416:GLN:HG3	3:A:469:HOH:O	2.04	0.58
1:A:256:LEU:HA	1:A:283:ALA:O	2.04	0.57
1:D:134:LEU:HD22	1:D:185:ILE:HG22	1.85	0.57
1:D:286:VAL:HG23	1:D:298:VAL:CG2	2.27	0.57
1:A:145:THR:CG2	1:A:146:LEU:N	2.67	0.57
1:C:114:PHE:O	1:C:118:VAL:HG23	2.04	0.57
1:B:230:GLY:H	1:B:233:SER:HB3	1.69	0.57
1:A:367:ASP:OD2	1:A:370:ARG:HD3	2.04	0.57
1:D:165:ARG:NH1	1:D:211:PHE:HA	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:416:GLN:HE21	1:C:416:GLN:HA	1.69	0.56
1:C:287:SER:HB3	1:C:325:ASP:O	2.05	0.56
1:C:357:LYS:CD	1:C:357:LYS:H	2.16	0.56
1:D:195:ASP:CG	1:D:198:THR:HG23	2.29	0.56
1:D:280:ASP:OD1	1:D:281:LYS:N	2.38	0.56
1:B:400:GLU:O	1:B:403:LEU:HB2	2.05	0.56
1:A:299:THR:HA	1:A:343:ASN:ND2	2.20	0.56
1:D:83:LEU:HD22	1:D:397:PRO:HG2	1.87	0.56
1:A:412:GLU:HG2	1:A:413:VAL:HG22	1.87	0.56
1:D:27:ARG:HB2	3:D:456:HOH:O	2.06	0.56
1:C:68:THR:HG22	1:C:69:THR:HG23	1.87	0.56
1:A:106:ARG:HA	1:A:109:LEU:HD13	1.86	0.55
1:A:74:ARG:NH2	1:A:98:GLU:OE1	2.31	0.55
1:B:98:GLU:C	1:B:99:THR:O	2.47	0.55
1:C:33:ILE:H	1:C:88:HIS:CD2	2.25	0.55
1:C:138:GLU:O	1:C:141:LYS:HG2	2.07	0.55
1:C:145:THR:HB	1:C:187:ASP:OD1	2.07	0.55
1:B:14:TYR:O	1:B:19:ILE:HB	2.07	0.55
1:B:106:ARG:HA	1:B:109:LEU:HD22	1.89	0.55
1:A:234:GLU:CD	1:D:292:TRP:HE1	2.13	0.55
1:C:261:HIS:HE1	3:C:455:HOH:O	1.88	0.55
1:A:309:LEU:O	1:A:313:ARG:HB2	2.07	0.54
1:D:145:THR:CG2	1:D:146:LEU:N	2.68	0.54
1:A:85:PRO:HD3	1:A:396:VAL:HG21	1.90	0.54
1:C:145:THR:HG22	1:C:146:LEU:N	2.12	0.54
1:C:198:THR:N	1:C:199:PRO:HD2	2.21	0.54
1:D:261:HIS:CE1	1:D:293:ASP:OD2	2.61	0.54
1:D:280:ASP:HA	1:D:319:ARG:NH1	2.22	0.54
1:C:245:LEU:HD22	1:D:245:LEU:HD23	1.85	0.54
1:C:402:TRP:CZ3	1:C:403:LEU:HD13	2.43	0.54
1:D:145:THR:HG22	1:D:146:LEU:HG	1.90	0.54
1:A:245:LEU:HD22	1:B:245:LEU:HD22	1.90	0.54
1:C:334:ILE:HG22	1:C:406:ILE:CD1	2.39	0.53
1:A:187:ASP:HB3	1:A:238:VAL:HG21	1.91	0.53
1:A:197:LEU:HD11	1:B:277:LEU:CD1	2.39	0.53
1:A:104:VAL:CG2	1:A:109:LEU:HD12	2.39	0.53
1:D:226:SER:O	1:D:261:HIS:HD2	1.90	0.53
1:C:413:VAL:O	1:C:414:LEU:C	2.50	0.53
1:D:280:ASP:HA	1:D:319:ARG:HH11	1.74	0.53
1:A:227:LYS:HE2	1:A:261:HIS:CE1	2.43	0.53
1:D:180:LEU:HD12	1:D:221:LEU:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:334:ILE:HG22	1:B:406:ILE:HD12	1.90	0.53
1:D:102:LYS:O	1:D:103:VAL:HG12	2.09	0.53
1:D:148:HIS:ND1	1:D:149:PRO:CD	2.69	0.53
1:C:359:LEU:HD21	1:C:374:VAL:HG22	1.90	0.52
1:D:261:HIS:HE1	1:D:293:ASP:OD2	1.93	0.52
1:D:110:GLU:H	1:D:113:HIS:CD2	2.16	0.52
1:D:8:GLU:OE1	1:D:11:LYS:HE2	2.10	0.52
1:A:168:GLY:O	1:A:171:PHE:HB2	2.10	0.52
1:C:261:HIS:CE1	1:C:293:ASP:OD2	2.62	0.52
1:C:400:GLU:O	1:C:403:LEU:HB2	2.10	0.52
1:B:234:GLU:OE1	1:C:292:TRP:NE1	2.42	0.52
1:B:361:GLU:O	1:B:365:THR:HB	2.09	0.52
1:C:148:HIS:CE1	1:C:149:PRO:HD2	2.45	0.52
1:A:303:GLU:O	1:A:307:ILE:HG13	2.10	0.52
1:D:196:ARG:HG3	1:D:196:ARG:HH11	1.75	0.52
1:A:197:LEU:HD11	1:B:277:LEU:HD13	1.92	0.51
1:A:77:LEU:O	1:A:81:LEU:HD22	2.09	0.51
1:A:413:VAL:O	1:A:414:LEU:C	2.53	0.51
1:C:229:PHE:HB2	1:C:234:GLU:HA	1.93	0.51
1:C:324:LEU:HD21	1:C:344:VAL:HG11	1.91	0.51
1:D:370:ARG:CG	1:D:370:ARG:NH1	2.64	0.51
1:A:110:GLU:N	1:A:113:HIS:HD2	2.00	0.51
1:A:410:GLU:CA	1:A:414:LEU:HB2	2.26	0.51
1:B:172:GLY:HA2	1:B:177:THR:O	2.11	0.51
1:B:180:LEU:HD11	1:B:223:ALA:HB2	1.93	0.51
1:A:158:ILE:HG23	1:A:210:ILE:HG13	1.91	0.51
1:A:285:HIS:HB3	1:A:325:ASP:HB2	1.92	0.51
1:A:161:CYS:HB2	1:A:210:ILE:HG12	1.92	0.50
1:B:33:ILE:H	1:B:88:HIS:HD2	1.59	0.50
1:D:21:VAL:HG13	3:D:481:HOH:O	2.10	0.50
1:D:134:LEU:CD2	1:D:145:THR:HG21	2.39	0.50
1:D:232:GLY:N	1:D:234:GLU:OE2	2.35	0.50
1:A:187:ASP:C	1:A:238:VAL:HG23	2.35	0.50
1:B:256:LEU:HA	1:B:283:ALA:O	2.11	0.50
1:B:143:GLY:O	1:B:187:ASP:HA	2.11	0.50
1:D:257:LEU:HD11	1:D:275:MET:HE2	1.93	0.50
1:C:145:THR:HG23	1:C:146:LEU:N	2.26	0.50
1:A:357:LYS:CD	1:A:357:LYS:H	2.24	0.49
1:A:359:LEU:HD21	1:A:374:VAL:HG22	1.94	0.49
1:B:103:VAL:O	1:B:103:VAL:HG13	2.11	0.49
1:A:39:GLN:HB3	1:A:337:TRP:NE1	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:HIS:CE1	1:A:131:ASN:HB2	2.47	0.49
1:C:300:PHE:HE1	1:C:347:ALA:HA	1.77	0.49
1:B:230:GLY:N	1:B:233:SER:HB3	2.27	0.49
1:D:99:THR:O	1:D:100:ASP:O	2.30	0.49
1:C:279:HIS:O	1:C:280:ASP:CG	2.56	0.49
1:A:400:GLU:HG3	1:C:377:GLU:HG3	1.94	0.49
1:B:85:PRO:O	1:B:88:HIS:HE1	1.95	0.49
1:A:42:ASP:OD2	1:A:333:ARG:NH2	2.45	0.49
1:A:370:ARG:O	1:A:374:VAL:HG13	2.13	0.49
1:D:410:GLU:HA	1:D:414:LEU:HB2	1.95	0.49
1:C:106:ARG:HA	1:C:109:LEU:HD22	1.95	0.48
1:A:66:LYS:HB3	1:A:333:ARG:CZ	2.43	0.48
1:B:224:VAL:HB	1:B:244:TYR:CD1	2.48	0.48
1:A:324:LEU:HD13	1:A:326:PHE:CE2	2.49	0.48
1:C:200:ARG:NH1	1:D:273:SER:OG	2.45	0.48
1:C:257:LEU:HD11	1:C:275:MET:CE	2.37	0.48
1:C:119:ARG:CB	1:C:119:ARG:HH11	2.27	0.48
1:B:31:VAL:HG11	1:B:317:LEU:HD22	1.96	0.48
1:C:95:ILE:CG2	1:C:133:THR:HG21	2.44	0.48
1:C:364:GLU:CD	1:D:202:ARG:HH22	2.21	0.48
1:B:43:VAL:HG12	1:B:43:VAL:O	2.14	0.47
1:B:111:PRO:HB3	1:B:170:TYR:CD2	2.49	0.47
1:A:321:LEU:N	1:A:321:LEU:CD1	2.78	0.47
1:B:13:GLU:OE2	1:D:399:LYS:NZ	2.43	0.47
1:B:39:GLN:HE22	1:B:328:ASP:H	1.62	0.47
1:B:402:TRP:CZ3	1:B:403:LEU:HD13	2.49	0.47
1:C:95:ILE:HG23	1:C:133:THR:HG21	1.96	0.47
1:D:290:VAL:O	1:D:291:ARG:HB2	2.15	0.47
1:D:292:TRP:O	1:D:294:SER:N	2.48	0.47
1:B:110:GLU:OE1	1:B:112:ARG:NH1	2.48	0.47
1:B:33:ILE:O	1:B:88:HIS:HB3	2.14	0.47
1:B:258:ASP:HB3	1:B:261:HIS:CG	2.50	0.47
1:B:412:GLU:O	1:B:416:GLN:HG3	2.15	0.47
1:D:29:LYS:HG3	1:D:394:MET:HE1	1.96	0.47
1:D:102:LYS:O	1:D:103:VAL:CG1	2.63	0.47
1:B:257:LEU:HD11	1:B:275:MET:HE2	1.97	0.46
1:B:280:ASP:OD1	1:B:281:LYS:N	2.48	0.46
1:C:21:VAL:HG12	1:C:353:LEU:HD22	1.97	0.46
1:C:36:HIS:HA	1:C:91:ASN:HB3	1.97	0.46
1:D:145:THR:HG23	1:D:157:TRP:CD1	2.50	0.46
1:A:296:HIS:HE1	1:D:234:GLU:HG3	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:THR:HG23	1:D:157:TRP:CG	2.50	0.46
1:D:377:GLU:HA	1:D:377:GLU:OE1	2.15	0.46
1:A:48:LEU:N	1:A:48:LEU:CD1	2.78	0.46
1:B:111:PRO:HB3	1:B:170:TYR:CG	2.51	0.46
1:D:84:ILE:HG22	1:D:88:HIS:CE1	2.51	0.46
1:A:300:PHE:HE1	1:A:347:ALA:HA	1.81	0.46
1:C:266:GLU:HG3	1:D:263:HIS:CE1	2.51	0.46
1:D:168:GLY:O	1:D:171:PHE:HB2	2.16	0.46
1:B:334:ILE:HG22	1:B:406:ILE:CD1	2.46	0.46
1:A:227:LYS:HA	1:A:261:HIS:HB3	1.97	0.45
1:B:9:ARG:HD2	1:D:6:GLN:NE2	2.31	0.45
1:B:309:LEU:O	1:B:313:ARG:HG3	2.17	0.45
1:C:84:ILE:HG21	1:C:341:THR:HG21	1.97	0.45
1:D:232:GLY:H	1:D:234:GLU:CD	2.21	0.45
1:B:33:ILE:N	1:B:88:HIS:HD2	2.13	0.45
1:A:301:ASP:HB2	1:C:302:ASP:OD2	2.16	0.45
1:A:417:ARG:H	1:A:417:ARG:HG3	1.51	0.45
1:A:148:HIS:ND1	1:A:149:PRO:CD	2.78	0.45
1:B:110:GLU:N	1:B:113:HIS:HD2	2.02	0.45
1:C:227:LYS:HA	1:C:261:HIS:HB3	1.98	0.45
1:A:9:ARG:NH1	1:C:3:MET:N	2.64	0.45
1:D:41:ASP:HB2	1:D:67:ALA:O	2.16	0.45
1:A:295:ASP:HB3	1:A:327:PHE:N	2.30	0.45
1:D:36:HIS:NE2	1:D:326:PHE:O	2.49	0.45
1:D:288:ARG:HB3	1:D:296:HIS:HB2	1.99	0.45
1:A:331:ILE:HG22	1:C:368:TYR:HB2	1.98	0.45
1:C:258:ASP:HB3	1:C:261:HIS:CG	2.52	0.45
1:C:309:LEU:O	1:C:313:ARG:HB3	2.17	0.45
1:D:134:LEU:HD23	1:D:145:THR:CG2	2.40	0.45
1:D:234:GLU:H	1:D:234:GLU:HG2	1.47	0.45
1:C:42:ASP:HA	1:C:327:PHE:CE2	2.52	0.45
1:C:299:THR:HA	1:C:343:ASN:ND2	2.32	0.45
1:A:106:ARG:NH1	1:A:133:THR:O	2.50	0.45
1:D:148:HIS:CG	1:D:149:PRO:HD2	2.50	0.45
1:C:36:HIS:NE2	1:C:326:PHE:O	2.48	0.44
1:D:224:VAL:HB	1:D:244:TYR:CD1	2.52	0.44
1:D:250:LYS:HA	1:D:250:LYS:HD2	1.77	0.44
1:D:110:GLU:N	1:D:113:HIS:HD2	2.05	0.44
1:A:129:ASP:HA	1:A:179:CYS:SG	2.58	0.44
1:C:413:VAL:O	1:C:415:LEU:N	2.50	0.44
1:D:282:LEU:O	1:D:320:VAL:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:30:GLN:H	1:C:30:GLN:HG2	1.69	0.44
1:C:200:ARG:HH21	1:C:240:SER:HB3	1.82	0.44
1:A:397:PRO:HB2	1:A:402:TRP:HB3	2.00	0.44
1:B:229:PHE:CD1	1:C:291:ARG:HD2	2.53	0.44
1:B:234:GLU:CD	1:C:292:TRP:HE1	2.26	0.44
1:D:70:PRO:HB2	1:D:74:ARG:NH2	2.33	0.44
1:A:357:LYS:H	1:A:357:LYS:HD3	1.82	0.44
1:B:33:ILE:H	1:B:88:HIS:CD2	2.36	0.44
1:C:288:ARG:HB3	1:C:296:HIS:HB2	2.00	0.43
1:B:295:ASP:HB3	1:B:327:PHE:H	1.82	0.43
1:A:271:LYS:HB3	1:A:275:MET:CE	2.47	0.43
1:B:26:GLU:OE1	1:B:26:GLU:HA	2.17	0.43
1:A:286:VAL:CG2	1:A:298:VAL:HG21	2.36	0.43
1:B:145:THR:HB	1:B:187:ASP:OD1	2.18	0.43
1:C:261:HIS:HE1	1:C:293:ASP:OD2	2.02	0.43
1:C:184:TRP:CZ2	1:C:227:LYS:HE3	2.54	0.43
1:D:227:LYS:HD3	1:D:229:PHE:O	2.19	0.43
1:B:370:ARG:HG3	1:B:370:ARG:NH1	2.11	0.43
1:D:376:GLU:HA	1:D:379:LYS:HD2	2.01	0.43
1:C:256:LEU:HD11	1:C:285:HIS:NE2	2.34	0.43
1:D:305:ARG:HG3	1:D:305:ARG:NH1	2.34	0.42
1:A:151:GLN:HG2	1:A:154:ARG:NH2	2.35	0.42
1:A:381:TYR:OH	1:C:400:GLU:OE1	2.34	0.42
1:B:275:MET:HG2	1:B:282:LEU:HD21	2.01	0.42
1:D:89:ARG:HH21	1:D:129:ASP:CG	2.27	0.42
1:C:100:ASP:HB3	1:C:101:GLY:H	1.66	0.42
1:A:280:ASP:HA	1:A:319:ARG:HH11	1.81	0.42
1:D:399:LYS:HB3	1:D:400:GLU:H	1.63	0.42
1:C:110:GLU:HA	3:C:449:HOH:O	2.20	0.42
1:C:334:ILE:CG2	1:C:406:ILE:HD13	2.50	0.42
1:A:42:ASP:HA	1:A:327:PHE:CE2	2.55	0.42
1:A:247:TYR:C	1:A:247:TYR:CD2	2.97	0.42
1:D:82:SER:HB2	1:D:83:LEU:HG	2.02	0.42
1:D:226:SER:O	1:D:261:HIS:CD2	2.71	0.42
1:A:282:LEU:HD12	1:A:320:VAL:HG22	2.01	0.42
1:B:36:HIS:ND1	1:B:36:HIS:N	2.67	0.42
1:B:103:VAL:O	1:B:103:VAL:CG1	2.67	0.42
1:B:6:GLN:HG2	1:B:7:PHE:H	1.84	0.42
1:B:36:HIS:NE2	1:B:326:PHE:O	2.53	0.42
1:B:100:ASP:HB3	1:B:102:LYS:HG3	2.01	0.42
1:C:21:VAL:O	1:C:25:LEU:HG	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:305:ARG:HG3	1:D:305:ARG:HH11	1.85	0.42
1:A:196:ARG:O	1:A:200:ARG:HD2	2.20	0.41
1:B:68:THR:HG22	1:B:69:THR:HG23	2.00	0.41
1:C:33:ILE:H	1:C:88:HIS:HD2	1.65	0.41
1:C:222:ASP:O	1:C:254:LEU:HB2	2.20	0.41
1:C:334:ILE:HG22	1:C:406:ILE:HD11	2.02	0.41
1:A:310:GLU:OE2	1:A:313:ARG:NH1	2.53	0.41
1:A:356:HIS:O	1:A:360:LYS:HG3	2.21	0.41
1:B:250:LYS:HD3	1:B:250:LYS:HA	1.64	0.41
1:C:367:ASP:CG	1:C:370:ARG:HG2	2.45	0.41
1:C:394:MET:HA	1:C:394:MET:CE	2.51	0.41
1:C:397:PRO:HG2	1:C:402:TRP:HA	2.03	0.41
1:A:104:VAL:HG22	1:A:109:LEU:HD12	2.01	0.41
1:B:182:ASN:HD21	1:B:225:GLU:HB2	1.84	0.41
1:C:11:LYS:HD3	1:C:21:VAL:CG2	2.51	0.41
1:A:357:LYS:HG2	1:A:358:GLN:OE1	2.20	0.41
1:C:304:LEU:HD21	1:C:344:VAL:HA	2.03	0.41
1:C:31:VAL:HA	1:C:32:PRO:HD3	1.68	0.41
1:C:377:GLU:OE1	1:C:377:GLU:HA	2.20	0.41
1:D:180:LEU:HD11	1:D:223:ALA:HB2	2.01	0.41
1:B:28:LEU:HD23	1:B:390:TYR:CZ	2.56	0.41
1:C:202:ARG:HH11	1:C:202:ARG:HD3	1.76	0.41
1:D:99:THR:O	1:D:100:ASP:C	2.62	0.41
1:A:148:HIS:O	1:A:154:ARG:HD3	2.20	0.41
1:A:182:ASN:HD21	1:A:225:GLU:HB2	1.85	0.41
1:A:251:ASN:O	1:A:253:LYS:HG2	2.20	0.41
1:A:324:LEU:HD21	1:A:344:VAL:HG11	2.02	0.41
1:B:288:ARG:HA	1:B:289:PRO:HD2	1.94	0.41
1:C:272:ILE:HG22	1:C:276:LEU:HD22	2.02	0.41
1:C:273:SER:OG	1:D:200:ARG:NH1	2.54	0.41
1:C:409:TYR:CD2	1:C:414:LEU:HG	2.56	0.41
1:D:33:ILE:H	1:D:88:HIS:CD2	2.25	0.41
1:A:224:VAL:HB	1:A:244:TYR:CD1	2.56	0.41
1:A:234:GLU:HB3	1:D:290:VAL:HB	2.02	0.41
1:A:293:ASP:CG	3:A:462:HOH:O	2.64	0.41
1:B:300:PHE:HE1	1:B:347:ALA:HA	1.86	0.41
1:B:340:GLY:O	1:B:344:VAL:HG12	2.21	0.41
1:C:33:ILE:O	1:C:88:HIS:HB3	2.21	0.41
1:D:231:ILE:CD1	1:D:292:TRP:CZ3	3.04	0.40
1:C:109:LEU:HD12	1:C:113:HIS:CD2	2.56	0.40
1:C:227:LYS:HG2	1:C:261:HIS:CG	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:PRO:HA	1:A:88:HIS:HD2	1.85	0.40
1:A:350:PHE:CD1	1:A:350:PHE:C	2.99	0.40
1:B:319:ARG:NH2	3:B:454:HOH:O	2.55	0.40
1:C:119:ARG:HH11	1:C:119:ARG:HB3	1.86	0.40
1:C:354:ILE:HG23	1:C:355:PRO:HD2	2.02	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	398/424 (94%)	359 (90%)	33 (8%)	6 (2%)	8	14
1	B	398/424 (94%)	367 (92%)	27 (7%)	4 (1%)	12	22
1	C	400/424 (94%)	370 (92%)	22 (6%)	8 (2%)	6	10
1	D	400/424 (94%)	365 (91%)	30 (8%)	5 (1%)	9	16
All	All	1596/1696 (94%)	1461 (92%)	112 (7%)	23 (1%)	9	15

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	48	LEU
1	B	100	ASP
1	B	280	ASP
1	C	412	GLU
1	D	100	ASP
1	A	103	VAL
1	C	100	ASP
1	C	112	ARG
1	D	364	GLU
1	A	29	LYS

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Mol	Chain	Res	Type
1	B	103	VAL
1	C	62	ASP
1	C	414	LEU
1	D	293	ASP
1	A	293	ASP
1	A	100	ASP
1	C	300	PHE
1	C	48	LEU
1	D	399	LYS
1	D	103	VAL
1	B	43	VAL
1	A	43	VAL
1	C	43	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	350/369 (95%)	301 (86%)	49 (14%)	3	5
1	B	350/369 (95%)	302 (86%)	48 (14%)	3	6
1	C	352/369 (95%)	303 (86%)	49 (14%)	3	5
1	D	352/369 (95%)	304 (86%)	48 (14%)	3	6
All	All	1404/1476 (95%)	1210 (86%)	194 (14%)	3	5

All (194) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	21	VAL
1	A	31	VAL
1	A	43	VAL
1	A	48	LEU
1	A	68	THR
1	A	75	MET
1	A	79	LYS

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Mol	Chain	Res	Type
1	A	81	LEU
1	A	87	LYS
1	A	89	ARG
1	A	93	HIS
1	A	103	VAL
1	A	104	VAL
1	A	131	ASN
1	A	141	LYS
1	A	145	THR
1	A	151	GLN
1	A	175	LEU
1	A	198	THR
1	A	209	GLN
1	A	210	ILE
1	A	214	GLU
1	A	233	SER
1	A	275	MET
1	A	276	LEU
1	A	277	LEU
1	A	280	ASP
1	A	281	LYS
1	A	282	LEU
1	A	284	LEU
1	A	310	GLU
1	A	312	VAL
1	A	313	ARG
1	A	321	LEU
1	A	338	THR
1	A	344	VAL
1	A	354	ILE
1	A	357	LYS
1	A	363	GLN
1	A	370	ARG
1	A	374	VAL
1	A	398	ILE
1	A	401	GLU
1	A	403	LEU
1	A	406	ILE
1	A	411	LYS
1	A	412	GLU
1	A	413	VAL
1	B	5	SER

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Mol	Chain	Res	Type
1	B	6	GLN
1	B	12	ILE
1	B	16	GLN
1	B	21	VAL
1	B	30	GLN
1	B	31	VAL
1	B	47	GLU
1	B	48	LEU
1	B	66	LYS
1	B	68	THR
1	B	79	LYS
1	B	81	LEU
1	B	87	LYS
1	B	104	VAL
1	B	109	LEU
1	B	115	GLU
1	B	141	LYS
1	B	145	THR
1	B	162	ILE
1	B	167	ILE
1	B	175	LEU
1	B	209	GLN
1	B	234	GLU
1	B	245	LEU
1	B	250	LYS
1	B	273	SER
1	B	276	LEU
1	B	277	LEU
1	B	281	LYS
1	B	282	LEU
1	B	290	VAL
1	B	321	LEU
1	B	334	ILE
1	B	338	THR
1	B	344	VAL
1	B	357	LYS
1	B	360	LYS
1	B	364	GLU
1	B	365	THR
1	B	370	ARG
1	B	374	VAL
1	B	394	MET

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Mol	Chain	Res	Type
1	B	398	ILE
1	B	401	GLU
1	B	403	LEU
1	B	404	LYS
1	B	412	GLU
1	C	3	MET
1	C	16	GLN
1	C	23	GLU
1	C	27	ARG
1	C	29	LYS
1	C	30	GLN
1	C	31	VAL
1	C	43	VAL
1	C	48	LEU
1	C	68	THR
1	C	81	LEU
1	C	87	LYS
1	C	100	ASP
1	C	103	VAL
1	C	104	VAL
1	C	109	LEU
1	C	116	LYS
1	C	119	ARG
1	C	139	LYS
1	C	145	THR
1	C	175	LEU
1	C	183	ILE
1	C	227	LYS
1	C	231	ILE
1	C	234	GLU
1	C	237	VAL
1	C	269	SER
1	C	275	MET
1	C	276	LEU
1	C	277	LEU
1	C	280	ASP
1	C	281	LYS
1	C	282	LEU
1	C	286	VAL
1	C	313	ARG
1	C	321	LEU
1	C	338	THR

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Mol	Chain	Res	Type
1	C	344	VAL
1	C	357	LYS
1	C	364	GLU
1	C	365	THR
1	C	374	VAL
1	C	394	MET
1	C	399	LYS
1	C	403	LEU
1	C	404	LYS
1	C	406	ILE
1	C	413	VAL
1	C	416	GLN
1	D	4	LYS
1	D	6	GLN
1	D	21	VAL
1	D	30	GLN
1	D	31	VAL
1	D	43	VAL
1	D	68	THR
1	D	75	MET
1	D	79	LYS
1	D	81	LEU
1	D	93	HIS
1	D	99	THR
1	D	100	ASP
1	D	102	LYS
1	D	109	LEU
1	D	116	LYS
1	D	131	ASN
1	D	145	THR
1	D	151	GLN
1	D	164	SER
1	D	175	LEU
1	D	198	THR
1	D	201	LYS
1	D	214	GLU
1	D	215	ILE
1	D	217	GLU
1	D	224	VAL
1	D	227	LYS
1	D	234	GLU
1	D	250	LYS

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Mol	Chain	Res	Type
1	D	275	MET
1	D	276	LEU
1	D	280	ASP
1	D	282	LEU
1	D	284	LEU
1	D	315	ASP
1	D	321	LEU
1	D	338	THR
1	D	344	VAL
1	D	361	GLU
1	D	364	GLU
1	D	365	THR
1	D	370	ARG
1	D	374	VAL
1	D	401	GLU
1	D	403	LEU
1	D	404	LYS
1	D	413	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	88	HIS
1	A	113	HIS
1	A	131	ASN
1	A	182	ASN
1	A	251	ASN
1	A	388	ASN
1	A	416	GLN
1	B	39	GLN
1	B	88	HIS
1	B	93	HIS
1	B	113	HIS
1	B	131	ASN
1	B	182	ASN
1	B	314	ASN
1	B	388	ASN
1	C	39	GLN
1	C	88	HIS
1	C	113	HIS
1	C	182	ASN

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Mol	Chain	Res	Type
1	C	209	GLN
1	C	251	ASN
1	C	261	HIS
1	C	279	HIS
1	C	314	ASN
1	C	388	ASN
1	C	416	GLN
1	D	6	GLN
1	D	30	GLN
1	D	88	HIS
1	D	108	GLN
1	D	113	HIS
1	D	131	ASN
1	D	182	ASN
1	D	251	ASN
1	D	261	HIS
1	D	388	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	402/424 (94%)	-0.37	5 (1%) 76 75	16, 28, 47, 73	0
1	B	402/424 (94%)	-0.40	1 (0%) 91 92	18, 30, 42, 62	0
1	C	404/424 (95%)	-0.35	6 (1%) 72 69	16, 31, 49, 66	0
1	D	404/424 (95%)	-0.30	5 (1%) 76 75	17, 31, 54, 71	0
All	All	1612/1696 (95%)	-0.35	17 (1%) 78 77	16, 30, 49, 73	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	48	LEU	4.5
1	A	48	LEU	3.8
1	C	61	GLY	3.3
1	D	398	ILE	3.0
1	C	3	MET	3.0
1	C	413	VAL	2.8
1	C	49	SER	2.8
1	A	414	LEU	2.8
1	B	48	LEU	2.8
1	C	414	LEU	2.7
1	A	49	SER	2.4
1	D	49	SER	2.4
1	D	3	MET	2.2
1	D	48	LEU	2.2
1	D	417	ARG	2.1
1	A	413	VAL	2.1
1	A	38	ALA	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($\text{\AA}^2$)	Q<0.9
2	MN	A	419	1/1	0.90	0.16	83,83,83,83	0

6.5 Other polymers [i](#)

There are no such residues in this entry.