



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:59 AM UTC

PDB ID : 6JAK / pdb_00006jak
Title : OtsA apo structure
Authors : Wang, S.; Zhao, Y.; Wang, D.; Liu, J.
Deposited on : 2019-01-24
Resolution : 2.41 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

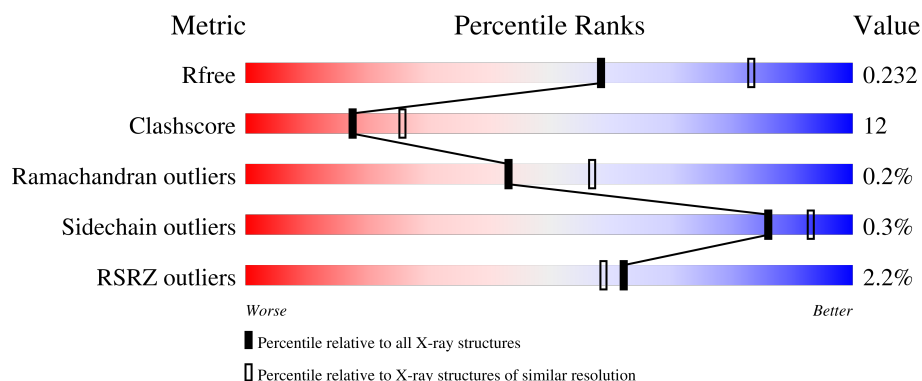
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	455	2% 82% 16% ..
1	B	455	3% 75% 22% ..
1	C	455	% 75% 22% ..
1	D	455	3% 75% 22% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14597 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 Trehalose-6-phosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	0	0
			3606	2319	626	654	7			
1	B	447	Total	C	N	O	S	0	0	0
			3583	2307	620	649	7			
1	C	447	Total	C	N	O	S	0	0	0
			3583	2307	620	649	7			
1	D	447	Total	C	N	O	S	0	0	0
			3583	2307	620	649	7			

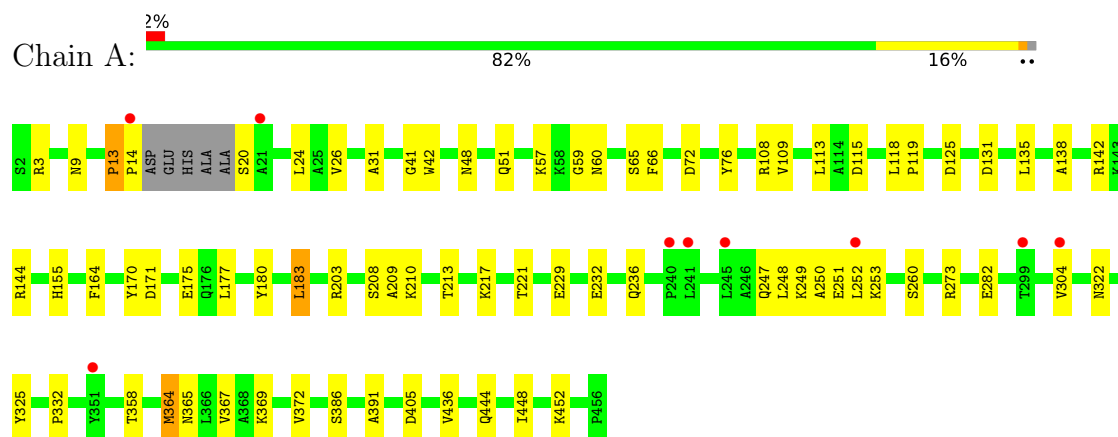
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	80	Total	O	0	0
			80	80		
2	B	53	Total	O	0	0
			53	53		
2	C	44	Total	O	0	0
			44	44		
2	D	65	Total	O	0	0
			65	65		

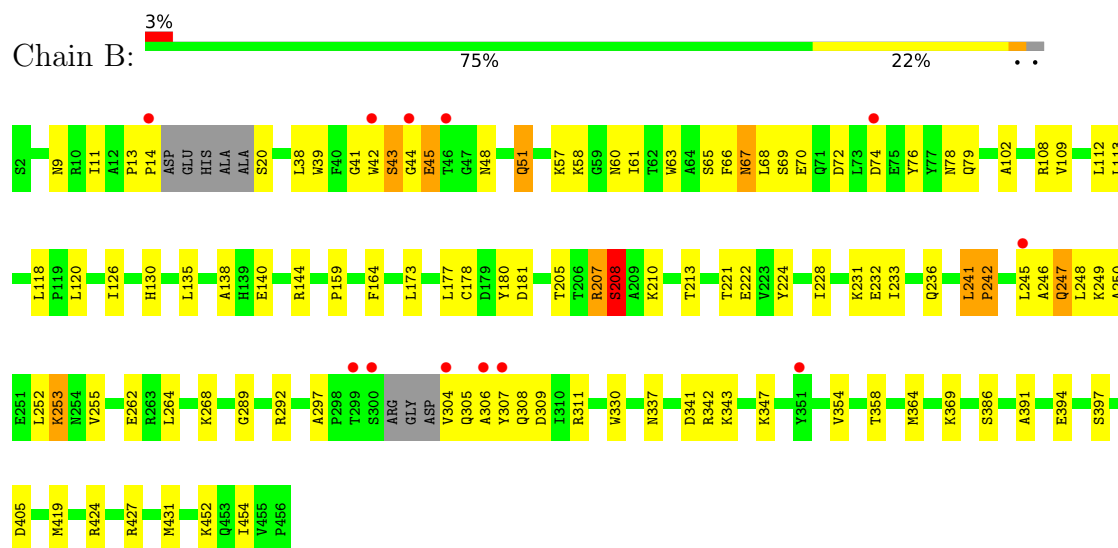
3 Residue-property plots

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.

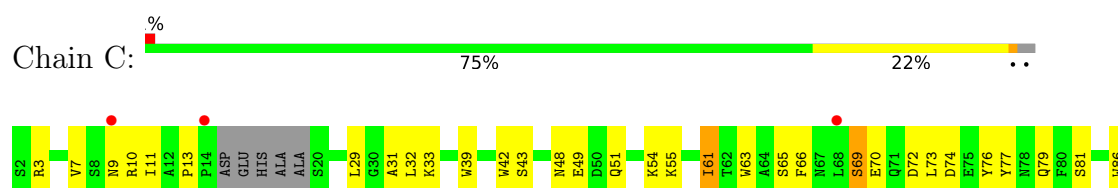
• Molecule 1: Trehalose-6-phosphate synthase

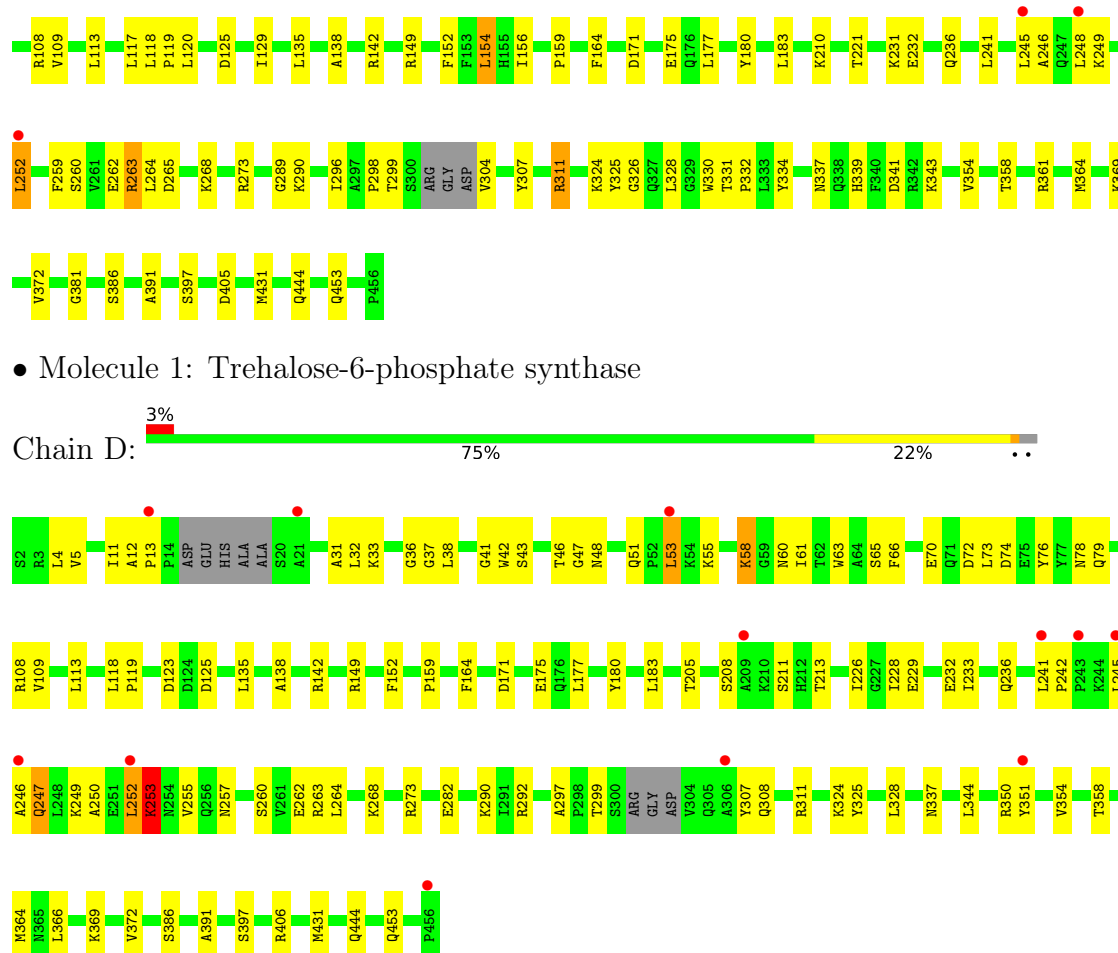


• Molecule 1: Trehalose-6-phosphate synthase



• Molecule 1: Trehalose-6-phosphate synthase





• Molecule 1: Trehalose-6-phosphate synthase

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.47Å 119.66Å 99.43Å 90.00° 91.88° 90.00°	Depositor
Resolution (Å)	29.69 – 2.41 29.69 – 2.41	Depositor EDS
% Data completeness (in resolution range)	95.3 (29.69-2.41) 95.3 (29.69-2.41)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.42Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.180 , 0.231 0.181 , 0.232	Depositor DCC
R_{free} test set	3736 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	50.7	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 29.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.043 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14597	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	0.56	0/3697	0.76	5/5026 (0.1%)
1	B	0.74	9/3673 (0.2%)	0.88	14/4993 (0.3%)
1	C	0.66	8/3673 (0.2%)	0.82	7/4993 (0.1%)
1	D	0.64	6/3673 (0.2%)	0.85	12/4993 (0.2%)
All	All	0.65	23/14716 (0.2%)	0.83	38/20005 (0.2%)

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	B	0	1
1	D	0	1
All	All	0	2

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	354	VAL	C-N	15.92	1.41	1.33
1	B	207	ARG	CZ-NH1	10.30	1.47	1.32
1	C	263	ARG	CZ-NH1	-10.25	1.18	1.32
1	C	354	VAL	C-N	9.89	1.38	1.33
1	B	67	ASN	CG-ND2	-9.28	1.13	1.33
1	C	263	ARG	NE-CZ	-7.93	1.24	1.33
1	B	242	PRO	N-CA	-7.59	1.36	1.46
1	C	263	ARG	CZ-NH2	-7.38	1.23	1.33
1	D	263	ARG	CZ-NH1	7.06	1.42	1.32
1	D	13	PRO	N-CA	-6.88	1.32	1.46
1	D	12	ALA	CA-CB	6.84	1.63	1.53
1	B	242	PRO	CA-C	6.71	1.55	1.51
1	C	311	ARG	CZ-NH2	-6.40	1.25	1.33
1	D	354	VAL	C-N	6.03	1.36	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	263	ARG	CD-NE	-5.87	1.38	1.46
1	B	207	ARG	CG-CD	-5.71	1.35	1.52
1	C	248	LEU	CG-CD1	-5.61	1.34	1.52
1	C	311	ARG	NE-CZ	-5.57	1.26	1.33
1	B	51	GLN	CD-NE2	-5.55	1.21	1.33
1	B	241	LEU	C-N	5.37	1.38	1.33
1	D	263	ARG	NE-CZ	5.29	1.38	1.33
1	B	45	GLU	CB-CG	-5.21	1.36	1.52
1	D	263	ARG	CZ-NH2	5.01	1.40	1.33

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	242	PRO	N-CA-CB	-12.97	95.93	103.19
1	D	58	LYS	CG-CD-CE	-11.25	85.42	111.30
1	C	290	LYS	CG-CD-CE	9.14	132.33	111.30
1	D	58	LYS	CD-CE-NZ	7.14	134.74	111.90
1	B	70	GLU	CB-CG-CD	-6.75	101.13	112.60
1	D	252	LEU	CA-C-N	-6.66	108.98	122.31
1	D	252	LEU	C-N-CA	-6.66	108.98	122.31
1	C	70	GLU	CB-CA-C	-6.61	99.82	110.79
1	B	207	ARG	CB-CA-C	-6.55	98.70	110.36
1	D	252	LEU	CB-CG-CD2	-6.36	91.63	110.70
1	D	263	ARG	NE-CZ-NH1	-6.30	115.20	121.50
1	D	13	PRO	N-CA-CB	-6.28	96.99	103.08
1	D	253	LYS	CG-CD-CE	6.20	125.55	111.30
1	D	13	PRO	CA-N-CD	6.14	120.60	112.00
1	B	68	LEU	N-CA-CB	-6.11	100.19	111.13
1	A	183	LEU	CB-CG-CD2	-5.89	93.02	110.70
1	D	13	PRO	CA-CB-CG	5.83	115.57	104.50
1	B	452	LYS	CA-CB-CG	5.57	125.24	114.10
1	B	452	LYS	CB-CA-C	-5.57	102.14	110.88
1	B	253	LYS	CG-CD-CE	-5.55	98.53	111.30
1	D	73	LEU	CA-CB-CG	-5.54	96.91	116.30
1	B	207	ARG	NE-CZ-NH1	-5.48	116.02	121.50
1	C	69	SER	CA-C-N	-5.40	113.05	120.28
1	C	69	SER	C-N-CA	-5.40	113.05	120.28
1	B	130	HIS	CA-C-N	-5.37	112.81	122.15
1	B	130	HIS	C-N-CA	-5.37	112.81	122.15
1	B	57	LYS	CA-CB-CG	-5.28	103.53	114.10
1	D	53	LEU	CB-CG-CD2	-5.24	94.98	110.70
1	B	207	ARG	NH1-CZ-NH2	5.23	126.09	119.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	LYS	CG-CD-CE	-5.22	99.30	111.30
1	C	61	ILE	CA-CB-CG1	5.21	119.26	110.40
1	A	364	MET	CA-C-N	-5.21	115.50	122.42
1	A	364	MET	C-N-CA	-5.21	115.50	122.42
1	A	304	VAL	CG1-CB-CG2	5.17	122.18	110.80
1	C	252	LEU	CA-CB-CG	5.17	134.38	116.30
1	B	208	SER	N-CA-C	-5.10	99.94	110.80
1	A	183	LEU	CD1-CG-CD2	-5.02	99.76	110.80
1	C	154	LEU	CA-CB-CG	5.00	133.81	116.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	247	GLN	Peptide
1	D	247	GLN	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	3606	0	3580	55	1
1	B	3583	0	3559	102	0
1	C	3583	0	3559	100	0
1	D	3583	0	3559	102	1
2	A	80	0	0	4	0
2	B	53	0	0	5	0
2	C	44	0	0	5	0
2	D	65	0	0	4	0
All	All	14597	0	14257	355	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

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:B:305:GLN:NE2	1:B:308:GLN:HE21	1.33	1.27
1:B:249:LYS:O	1:B:253:LYS:NZ	1.68	1.24
1:B:305:GLN:CD	1:B:308:GLN:NE2	1.98	1.21
1:C:263:ARG:HH21	1:C:298:PRO:HG2	1.07	1.16
1:D:252:LEU:HD22	1:D:255:VAL:CG1	1.77	1.14
1:D:252:LEU:HB2	1:D:255:VAL:HG12	1.29	1.14
1:B:305:GLN:CD	1:B:308:GLN:HE21	1.57	1.12
1:D:252:LEU:CD2	1:D:255:VAL:HG11	1.79	1.11
1:C:263:ARG:NH2	1:C:298:PRO:O	1.88	1.04
1:C:10:ARG:HH21	1:C:42:TRP:HZ3	1.02	0.98
1:C:263:ARG:HH21	1:C:298:PRO:CG	1.77	0.98
1:B:305:GLN:OE1	1:B:308:GLN:NE2	1.97	0.97
1:B:264:LEU:HD21	1:B:297:ALA:HB1	1.45	0.96
1:D:252:LEU:HD22	1:D:255:VAL:HG11	0.98	0.95
1:A:72:ASP:OD1	1:A:108:ARG:NH1	2.01	0.94
1:B:419:MET:SD	2:B:551:HOH:O	2.25	0.94
1:C:72:ASP:OD1	1:C:108:ARG:NH1	1.99	0.94
1:B:304:VAL:HG12	1:B:306:ALA:H	1.34	0.92
1:D:444:GLN:NE2	2:D:502:HOH:O	1.99	0.92
1:B:305:GLN:NE2	1:B:308:GLN:NE2	2.15	0.92
1:D:252:LEU:HB2	1:D:255:VAL:CG1	2.01	0.90
1:D:247:GLN:OE1	1:D:250:ALA:N	2.07	0.88
1:A:405:ASP:OD1	2:A:501:HOH:O	1.93	0.86
1:D:125:ASP:O	2:D:501:HOH:O	1.94	0.84
1:B:45:GLU:OE2	1:B:67:ASN:ND2	2.10	0.84
1:C:263:ARG:NH2	1:C:298:PRO:C	2.36	0.84
1:D:253:LYS:HA	1:D:253:LYS:HE2	1.58	0.84
1:D:48:ASN:O	1:D:51:GLN:NE2	2.10	0.83
1:D:249:LYS:HG3	1:D:351:TYR:CZ	2.14	0.82
1:B:205:THR:HG1	1:B:213:THR:HG1	1.22	0.82
1:D:46:THR:HG21	1:D:70:GLU:HG3	1.61	0.82
1:D:308:GLN:OE1	1:D:311:ARG:NH1	2.13	0.82
1:B:72:ASP:OD1	1:B:108:ARG:NH1	2.12	0.81
1:B:247:GLN:HB2	1:B:250:ALA:HB3	1.64	0.80
1:C:311:ARG:NH2	1:C:337:ASN:HD22	1.79	0.80
1:C:10:ARG:NH2	1:C:42:TRP:HZ3	1.80	0.80
1:D:72:ASP:OD1	1:D:108:ARG:NH1	2.15	0.79
1:C:263:ARG:NH2	1:C:298:PRO:HG2	1.93	0.79
1:B:364:MET:HE3	1:B:391:ALA:HA	1.63	0.79
1:C:263:ARG:HH22	1:C:298:PRO:C	1.90	0.79
1:D:252:LEU:HD21	1:D:292:ARG:NH2	1.97	0.78
1:D:33:LYS:HG2	1:D:61:ILE:HD11	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ARG:NH2	1:A:125:ASP:OD2	2.17	0.75
1:A:229:GLU:HG2	1:A:232:GLU:HB2	1.68	0.75
1:B:419:MET:HE2	1:B:424:ARG:HG2	1.68	0.74
1:B:307:TYR:OH	2:B:501:HOH:O	2.05	0.74
1:D:46:THR:HG21	1:D:70:GLU:CG	2.16	0.74
1:C:311:ARG:HH21	1:C:337:ASN:HD22	1.33	0.73
1:C:364:MET:HE2	1:C:369:LYS:HZ1	1.52	0.73
1:B:178:CYS:O	2:B:502:HOH:O	2.07	0.72
1:D:364:MET:HE2	1:D:369:LYS:NZ	2.04	0.72
1:D:252:LEU:CB	1:D:255:VAL:HG12	2.13	0.72
1:B:51:GLN:CD	1:B:51:GLN:H	1.96	0.72
1:C:364:MET:HE3	1:C:391:ALA:HA	1.71	0.72
1:D:252:LEU:HD21	1:D:292:ARG:CZ	2.21	0.71
1:D:364:MET:HE3	1:D:391:ALA:HA	1.72	0.71
1:B:241:LEU:HD12	1:B:241:LEU:H	1.55	0.70
1:D:46:THR:CG2	1:D:70:GLU:HG3	2.21	0.70
1:D:11:ILE:HD11	1:D:65:SER:HB2	1.73	0.70
1:D:242:PRO:O	1:D:246:ALA:N	2.24	0.70
1:D:241:LEU:HB2	1:D:246:ALA:HA	1.72	0.70
1:B:241:LEU:HB2	1:B:246:ALA:HA	1.74	0.69
1:A:13:PRO:HG2	1:A:14:PRO:HD3	1.74	0.69
1:A:364:MET:HE2	1:A:369:LYS:NZ	2.08	0.69
1:B:69:SER:OG	1:B:72:ASP:OD2	2.06	0.68
1:C:42:TRP:HB2	1:C:66:PHE:CE1	2.29	0.68
1:D:53:LEU:CD2	1:D:66:PHE:HB3	2.23	0.68
1:B:135:LEU:HD22	1:B:177:LEU:HD21	1.74	0.68
1:A:131:ASP:OD2	2:A:502:HOH:O	2.12	0.68
1:C:11:ILE:HD11	1:C:65:SER:HB3	1.74	0.68
1:C:304:VAL:N	2:C:506:HOH:O	2.27	0.67
1:B:9:ASN:O	1:B:42:TRP:HB3	1.94	0.67
1:B:250:ALA:HA	1:B:253:LYS:CE	2.24	0.67
1:B:311:ARG:HH21	1:B:337:ASN:HD21	1.43	0.67
1:D:253:LYS:HE3	1:D:351:TYR:OH	1.95	0.67
1:B:364:MET:HE2	1:B:369:LYS:NZ	2.08	0.67
1:C:381:GLY:O	2:C:502:HOH:O	2.12	0.66
1:A:208:SER:OG	1:A:209:ALA:N	2.29	0.65
1:B:13:PRO:HG2	1:B:14:PRO:HD3	1.77	0.65
1:C:154:LEU:HD13	1:C:156:ILE:O	1.96	0.65
1:C:405:ASP:OD1	2:C:503:HOH:O	2.15	0.65
1:C:74:ASP:OD2	1:C:79:GLN:NE2	2.29	0.65
1:B:140:GLU:O	1:B:144:ARG:HD3	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:241:LEU:HB2	1:C:246:ALA:HB2	1.79	0.64
1:C:328:LEU:HD12	1:D:245:LEU:HD22	1.79	0.64
1:B:305:GLN:HE22	1:B:308:GLN:HE21	1.42	0.64
1:D:299:THR:HA	1:D:307:TYR:OH	1.98	0.64
1:D:31:ALA:HB2	1:D:444:GLN:HG3	1.80	0.63
1:A:135:LEU:HD22	1:A:177:LEU:HD21	1.79	0.63
1:D:208:SER:HB2	1:D:211:SER:HB3	1.80	0.63
1:B:43:SER:HB3	1:B:45:GLU:OE2	1.99	0.63
1:B:252:LEU:HD22	1:B:255:VAL:HG21	1.81	0.63
1:B:364:MET:CE	1:B:391:ALA:HA	2.28	0.62
1:C:364:MET:HE2	1:C:369:LYS:CE	2.30	0.62
1:A:249:LYS:HA	1:A:252:LEU:HG	1.80	0.62
1:C:262:GLU:CG	1:C:268:LYS:HD2	2.31	0.61
1:C:361:ARG:NH1	2:C:501:HOH:O	2.07	0.61
1:C:263:ARG:NH1	1:C:307:TYR:HE1	1.99	0.61
1:B:245:LEU:HA	1:B:248:LEU:CD2	2.30	0.61
1:B:242:PRO:O	1:B:246:ALA:N	2.33	0.61
1:C:54:LYS:HD2	1:C:54:LYS:N	2.16	0.61
1:D:252:LEU:HD13	1:D:255:VAL:CG1	2.30	0.61
1:B:210:LYS:NZ	1:B:222:GLU:OE1	2.28	0.61
1:C:263:ARG:CZ	1:C:298:PRO:O	2.49	0.61
1:D:282:GLU:OE1	1:D:325:TYR:OH	2.07	0.60
1:A:364:MET:CE	1:A:391:ALA:HA	2.32	0.60
1:C:29:LEU:HB3	1:C:33:LYS:HE3	1.84	0.60
1:C:77:TYR:HA	1:C:81:SER:OG	2.02	0.60
1:D:42:TRP:HB2	1:D:66:PHE:CE1	2.36	0.60
1:C:358:THR:HA	1:C:386:SER:HB2	1.83	0.59
1:B:311:ARG:HH21	1:B:337:ASN:ND2	2.01	0.59
1:C:9:ASN:HD22	1:C:10:ARG:NH1	2.01	0.59
1:C:364:MET:HE2	1:C:369:LYS:NZ	2.17	0.59
1:B:210:LYS:HB3	1:B:221:THR:O	2.03	0.59
1:C:210:LYS:HB3	1:C:221:THR:O	2.03	0.59
1:B:308:GLN:N	1:B:308:GLN:OE1	2.36	0.59
1:A:364:MET:HE2	1:A:369:LYS:CE	2.33	0.58
1:D:228:ILE:HD12	1:D:233:ILE:HD13	1.86	0.58
1:B:311:ARG:NH2	1:B:337:ASN:HD21	2.01	0.58
1:C:263:ARG:NH2	1:C:298:PRO:CG	2.57	0.58
1:D:262:GLU:OE2	1:D:268:LYS:HD2	2.04	0.58
1:C:311:ARG:HD2	1:C:337:ASN:ND2	2.19	0.58
1:D:255:VAL:HG23	1:D:290:LYS:C	2.28	0.58
1:C:364:MET:CE	1:C:391:ALA:HA	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:364:MET:HE2	1:B:369:LYS:HZ1	1.66	0.57
1:A:138:ALA:HB2	1:A:180:TYR:CE1	2.40	0.57
1:C:263:ARG:NH1	1:C:298:PRO:O	2.38	0.57
1:C:311:ARG:NH2	1:C:337:ASN:ND2	2.49	0.56
1:D:43:SER:HB3	1:D:65:SER:OG	2.05	0.56
1:D:358:THR:HA	1:D:386:SER:HB2	1.86	0.56
1:B:364:MET:HE3	1:B:391:ALA:CA	2.35	0.56
1:C:397:SER:OG	1:C:431:MET:HG3	2.05	0.56
1:B:247:GLN:HB2	1:B:250:ALA:CB	2.33	0.56
1:A:57:LYS:NZ	1:A:59:GLY:O	2.28	0.56
1:C:299:THR:HG23	1:C:307:TYR:OH	2.06	0.55
1:B:250:ALA:HA	1:B:253:LYS:HE2	1.87	0.55
1:D:74:ASP:OD1	1:D:79:GLN:NE2	2.39	0.55
1:D:241:LEU:H	1:D:241:LEU:HD12	1.71	0.55
1:A:210:LYS:HB3	1:A:221:THR:O	2.07	0.55
1:C:29:LEU:O	1:C:33:LYS:HG3	2.07	0.55
1:C:117:LEU:HD12	1:C:120:LEU:HD12	1.87	0.55
1:B:252:LEU:HD11	1:B:292:ARG:HH12	1.72	0.55
1:D:252:LEU:CG	1:D:255:VAL:CG1	2.85	0.55
1:D:364:MET:HE2	1:D:369:LYS:HZ1	1.71	0.55
1:A:171:ASP:O	1:A:175:GLU:HG3	2.06	0.55
1:A:232:GLU:O	1:A:236:GLN:HG3	2.06	0.55
1:B:250:ALA:HA	1:B:253:LYS:NZ	2.22	0.54
1:B:262:GLU:CG	1:B:268:LYS:HD2	2.38	0.54
1:C:311:ARG:NH2	1:C:337:ASN:HB2	2.22	0.54
1:D:31:ALA:HB2	1:D:444:GLN:CG	2.37	0.54
1:D:138:ALA:HB2	1:D:180:TYR:CE1	2.42	0.54
1:B:343:LYS:HD2	1:B:347:LYS:HZ3	1.71	0.54
1:D:252:LEU:CD2	1:D:255:VAL:CG1	2.59	0.54
1:D:232:GLU:O	1:D:236:GLN:HG3	2.07	0.53
1:C:328:LEU:HD12	1:D:245:LEU:CD2	2.38	0.53
1:A:60:ASN:ND2	2:A:517:HOH:O	2.40	0.53
1:C:32:LEU:HB3	1:C:61:ILE:HD13	1.90	0.53
1:C:324:LYS:HD3	1:C:325:TYR:CE2	2.44	0.53
1:D:205:THR:OG1	1:D:213:THR:OG1	2.26	0.53
1:B:138:ALA:HB2	1:B:180:TYR:CE1	2.44	0.52
1:C:249:LYS:HD3	1:C:252:LEU:HD12	1.90	0.52
1:C:364:MET:HE3	1:C:391:ALA:CA	2.39	0.52
1:D:252:LEU:CB	1:D:255:VAL:CG1	2.80	0.52
1:C:311:ARG:NH2	1:C:337:ASN:CB	2.73	0.52
1:D:135:LEU:HD22	1:D:177:LEU:HD21	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:GLN:OE1	1:B:305:GLN:O	2.28	0.52
1:D:53:LEU:HD21	1:D:66:PHE:HB3	1.92	0.52
1:D:241:LEU:CB	1:D:246:ALA:HA	2.39	0.52
1:A:31:ALA:HB1	1:A:448:ILE:HD11	1.92	0.52
1:D:253:LYS:CE	1:D:351:TYR:OH	2.57	0.52
1:D:311:ARG:HH21	1:D:337:ASN:HD21	1.56	0.52
1:D:247:GLN:NE2	1:D:249:LYS:HB3	2.25	0.51
1:B:419:MET:HE3	1:B:424:ARG:HA	1.93	0.51
1:D:72:ASP:O	1:D:76:TYR:HB3	2.11	0.51
1:B:419:MET:HE1	1:B:427:ARG:HD2	1.91	0.51
1:B:241:LEU:HD13	1:B:246:ALA:HA	1.92	0.51
1:C:31:ALA:HB2	1:C:444:GLN:HG3	1.92	0.51
1:B:11:ILE:HD11	1:B:65:SER:HB3	1.92	0.51
1:C:11:ILE:HG22	1:C:39:TRP:NE1	2.26	0.51
1:B:343:LYS:HD2	1:B:343:LYS:C	2.36	0.51
1:C:263:ARG:NH2	1:C:298:PRO:CA	2.74	0.51
1:C:31:ALA:HB2	1:C:444:GLN:CG	2.42	0.50
1:D:118:LEU:HB3	1:D:119:PRO:HD3	1.93	0.50
1:D:397:SER:OG	1:D:431:MET:HG3	2.10	0.50
1:A:20:SER:HB3	1:A:26:VAL:HG22	1.92	0.50
1:A:109:VAL:O	1:A:113:LEU:HG	2.11	0.50
1:D:324:LYS:HD3	1:D:325:TYR:CE2	2.46	0.50
1:D:364:MET:CE	1:D:391:ALA:HA	2.41	0.50
1:C:54:LYS:HD2	1:C:54:LYS:H	1.76	0.49
1:C:73:LEU:HD12	1:C:77:TYR:HB3	1.94	0.49
1:C:289:GLY:HA2	1:C:330:TRP:CG	2.47	0.49
1:B:252:LEU:HD11	1:B:292:ARG:NH1	2.26	0.49
1:D:252:LEU:CD2	1:D:292:ARG:NH2	2.72	0.49
1:A:138:ALA:O	1:A:142:ARG:HG2	2.11	0.49
1:C:135:LEU:HD22	1:C:177:LEU:HD21	1.95	0.49
1:C:231:LYS:HG3	1:C:232:GLU:N	2.28	0.49
1:C:72:ASP:O	1:C:76:TYR:HB3	2.13	0.49
1:C:138:ALA:HB2	1:C:180:TYR:CE1	2.48	0.49
1:C:138:ALA:O	1:C:142:ARG:HG2	2.13	0.49
1:D:249:LYS:HG3	1:D:351:TYR:OH	2.11	0.49
1:A:364:MET:HE3	1:A:391:ALA:HB2	1.94	0.49
1:B:74:ASP:OD1	1:B:79:GLN:NE2	2.46	0.49
1:C:245:LEU:HD23	1:D:328:LEU:HD12	1.93	0.49
1:D:299:THR:HG23	1:D:337:ASN:OD1	2.12	0.49
1:A:48:ASN:HB3	1:A:51:GLN:HG2	1.94	0.49
1:C:152:PHE:HB3	1:C:183:LEU:HD23	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:299:THR:H	1:C:339:HIS:CD2	2.31	0.49
1:C:326:GLY:HA2	1:C:332:PRO:HD3	1.93	0.48
1:D:229:GLU:CD	1:D:232:GLU:HG3	2.37	0.48
1:D:5:VAL:HG13	1:D:38:LEU:HD12	1.93	0.48
1:A:260:SER:OG	1:A:273:ARG:NH1	2.46	0.48
1:B:72:ASP:HB3	1:B:109:VAL:CG2	2.43	0.48
1:D:171:ASP:O	1:D:175:GLU:HG3	2.13	0.48
1:D:252:LEU:CG	1:D:255:VAL:HG12	2.43	0.48
1:B:51:GLN:CD	1:B:51:GLN:N	2.68	0.48
1:C:149:ARG:NH1	1:C:453:GLN:O	2.36	0.48
1:A:364:MET:HE3	1:A:391:ALA:HA	1.94	0.48
1:B:289:GLY:HA2	1:B:330:TRP:CG	2.48	0.48
1:C:43:SER:OG	1:C:65:SER:HB2	2.14	0.48
1:B:43:SER:OG	1:B:65:SER:HB2	2.14	0.48
1:A:9:ASN:O	1:A:42:TRP:HB3	2.13	0.48
1:C:48:ASN:HB3	1:C:51:GLN:CD	2.39	0.48
1:C:117:LEU:O	1:C:117:LEU:HG	2.14	0.48
1:C:311:ARG:HH22	1:C:337:ASN:HB2	1.79	0.47
1:C:444:GLN:NE2	2:C:515:HOH:O	2.48	0.47
1:C:262:GLU:HG3	1:C:268:LYS:HD2	1.96	0.47
1:D:364:MET:HE3	1:D:391:ALA:CA	2.41	0.47
1:A:66:PHE:CZ	1:A:113:LEU:HD22	2.50	0.47
1:B:42:TRP:CD1	1:B:44:GLY:H	2.31	0.47
1:A:444:GLN:O	1:A:448:ILE:HD12	2.15	0.47
1:C:55:LYS:HA	1:C:63:TRP:O	2.15	0.47
1:B:11:ILE:HG23	1:B:41:GLY:HA3	1.96	0.47
1:B:228:ILE:HD12	1:B:233:ILE:HD12	1.95	0.47
1:C:249:LYS:NZ	1:C:252:LEU:O	2.35	0.47
1:C:263:ARG:HH12	1:C:307:TYR:HE1	1.62	0.47
1:D:260:SER:OG	1:D:273:ARG:NH1	2.48	0.47
1:B:72:ASP:O	1:B:76:TYR:HB3	2.14	0.47
1:B:358:THR:HA	1:B:386:SER:HB2	1.97	0.46
1:D:109:VAL:O	1:D:113:LEU:HG	2.15	0.46
1:A:364:MET:HE3	1:A:391:ALA:CB	2.45	0.46
1:C:10:ARG:NH2	1:C:77:TYR:CE2	2.70	0.46
1:B:181:ASP:HB3	1:B:454:ILE:HG21	1.96	0.46
1:B:245:LEU:HA	1:B:248:LEU:HD23	1.97	0.46
1:B:231:LYS:HA	1:B:231:LYS:HD3	1.72	0.46
1:D:11:ILE:HG23	1:D:41:GLY:HA3	1.98	0.46
1:D:55:LYS:HA	1:D:63:TRP:O	2.15	0.46
1:D:226:ILE:HG23	1:D:366:LEU:HD21	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:LEU:HB3	1:A:119:PRO:HD3	1.97	0.46
1:A:364:MET:HE1	1:A:391:ALA:HA	1.98	0.46
1:D:46:THR:HG22	1:D:47:GLY:N	2.31	0.46
1:C:262:GLU:HG2	1:C:268:LYS:HD2	1.98	0.46
1:D:262:GLU:CD	1:D:273:ARG:HH22	2.23	0.45
1:B:48:ASN:HB2	1:B:51:GLN:NE2	2.31	0.45
1:A:41:GLY:O	1:A:65:SER:HA	2.16	0.45
1:B:207:ARG:C	1:B:208:SER:O	2.54	0.45
1:D:33:LYS:NZ	1:D:60:ASN:HB2	2.31	0.45
1:D:4:LEU:O	1:D:37:GLY:HA3	2.16	0.45
1:C:66:PHE:CZ	1:C:113:LEU:HD22	2.51	0.45
1:D:123:ASP:OD1	1:D:123:ASP:N	2.49	0.45
1:B:79:GLN:OE1	1:B:102:ALA:HB2	2.17	0.45
1:B:264:LEU:HD21	1:B:297:ALA:CB	2.32	0.45
1:B:308:GLN:OE1	1:B:309:ASP:N	2.46	0.45
1:B:74:ASP:O	1:B:79:GLN:N	2.49	0.45
1:B:232:GLU:O	1:B:236:GLN:HG3	2.17	0.45
1:D:364:MET:HE2	1:D:369:LYS:CE	2.46	0.45
1:D:247:GLN:OE1	1:D:249:LYS:N	2.50	0.45
1:A:448:ILE:HG22	1:A:452:LYS:HE2	1.99	0.44
1:A:72:ASP:O	1:A:76:TYR:HB3	2.16	0.44
1:C:3:ARG:HB3	1:C:125:ASP:OD1	2.17	0.44
1:C:331:THR:HG23	1:C:334:TYR:CE1	2.52	0.44
1:C:118:LEU:HB3	1:C:119:PRO:HD3	2.00	0.44
1:D:350:ARG:HH11	1:D:350:ARG:HD2	1.63	0.44
1:A:247:GLN:HG3	1:A:248:LEU:HD13	2.00	0.44
1:D:58:LYS:HB2	1:D:63:TRP:CZ3	2.53	0.44
1:A:138:ALA:HB2	1:A:180:TYR:CD1	2.53	0.44
1:A:364:MET:HE3	1:A:391:ALA:CA	2.48	0.44
1:B:38:LEU:HD21	1:B:120:LEU:HD22	2.00	0.44
1:C:232:GLU:O	1:C:236:GLN:HG3	2.18	0.44
1:C:7:VAL:HB	1:C:129:ILE:HD13	2.00	0.44
1:C:54:LYS:NZ	1:C:66:PHE:HA	2.32	0.44
1:A:213:THR:HA	1:A:217:LYS:O	2.18	0.44
1:B:247:GLN:C	1:B:250:ALA:H	2.26	0.44
1:D:58:LYS:HB2	1:D:63:TRP:HZ3	1.83	0.44
1:A:248:LEU:HD12	1:A:251:GLU:HB2	1.98	0.44
1:B:419:MET:CE	1:B:424:ARG:HA	2.48	0.43
1:C:109:VAL:O	1:C:113:LEU:HG	2.18	0.43
1:D:48:ASN:HB3	1:D:51:GLN:NE2	2.33	0.43
1:A:250:ALA:HA	1:A:253:LYS:HE2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:46:THR:HG21	1:D:70:GLU:HG2	1.97	0.43
1:D:242:PRO:HG3	1:D:344:LEU:HD13	1.99	0.43
1:A:180:TYR:HB2	1:A:183:LEU:HD11	2.00	0.43
1:B:58:LYS:HB3	1:B:63:TRP:CZ3	2.53	0.43
1:B:405:ASP:OD1	2:B:503:HOH:O	2.21	0.43
1:D:53:LEU:HD23	1:D:66:PHE:HB3	1.97	0.43
1:A:365:ASN:OD1	1:A:367:VAL:HB	2.19	0.43
1:B:135:LEU:HD12	1:B:173:LEU:HD22	2.01	0.43
1:B:222:GLU:HB3	1:B:224:TYR:CE1	2.54	0.43
1:C:49:GLU:OE1	1:C:69:SER:HB3	2.18	0.43
1:B:66:PHE:CZ	1:B:113:LEU:HD22	2.53	0.43
1:C:260:SER:OG	1:C:273:ARG:NH1	2.52	0.43
1:D:138:ALA:O	1:D:142:ARG:HG2	2.19	0.43
1:A:252:LEU:O	1:A:252:LEU:HD12	2.19	0.42
1:A:436:VAL:O	2:A:503:HOH:O	2.22	0.42
1:D:32:LEU:O	1:D:36:GLY:N	2.52	0.42
1:A:322:ASN:OD1	1:A:332:PRO:HD2	2.20	0.42
1:B:13:PRO:CG	1:B:14:PRO:HD3	2.47	0.42
1:B:58:LYS:HB3	1:B:63:TRP:HZ3	1.85	0.42
1:C:263:ARG:NH1	1:C:307:TYR:CE1	2.83	0.42
1:A:13:PRO:CG	1:A:14:PRO:HD3	2.44	0.42
1:B:126:ILE:O	1:B:126:ILE:HG13	2.20	0.42
1:D:252:LEU:CD1	1:D:255:VAL:CG1	2.97	0.42
1:D:135:LEU:CD2	1:D:177:LEU:HD21	2.50	0.42
1:B:159:PRO:HG2	1:B:164:PHE:HB2	2.01	0.42
1:C:341:ASP:OD1	1:C:343:LYS:HG2	2.19	0.42
1:A:282:GLU:OE2	1:A:325:TYR:OH	2.35	0.42
1:D:149:ARG:NH1	1:D:453:GLN:O	2.51	0.42
1:B:74:ASP:HA	1:B:78:ASN:HB2	2.02	0.42
1:B:11:ILE:HG22	1:B:39:TRP:CD1	2.55	0.41
1:B:253:LYS:HZ3	1:B:253:LYS:HG2	1.77	0.41
1:B:419:MET:HE2	1:B:424:ARG:CG	2.45	0.41
1:D:264:LEU:HD21	1:D:297:ALA:HB1	2.02	0.41
1:A:31:ALA:HB1	1:A:448:ILE:CD1	2.49	0.41
1:A:358:THR:HA	1:A:386:SER:HB2	2.02	0.41
1:B:228:ILE:HD12	1:B:233:ILE:CD1	2.50	0.41
1:B:341:ASP:OD2	1:B:342:ARG:N	2.53	0.41
1:B:51:GLN:H	1:B:51:GLN:NE2	2.17	0.41
1:D:74:ASP:HA	1:D:78:ASN:HB2	2.03	0.41
1:D:152:PHE:HB3	1:D:183:LEU:HD23	2.01	0.41
1:C:86:TRP:CH2	1:C:265:ASP:OD1	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:249:LYS:HD3	1:C:249:LYS:HA	1.85	0.41
1:D:257:ASN:ND2	2:D:504:HOH:O	2.09	0.41
1:B:60:ASN:C	1:B:61:ILE:HG13	2.46	0.41
1:B:118:LEU:HA	1:B:118:LEU:HD12	1.85	0.41
1:C:72:ASP:CG	1:C:108:ARG:NH1	2.77	0.41
1:A:364:MET:HE2	1:A:369:LYS:HZ1	1.83	0.41
1:C:259:PHE:HZ	1:C:296:ILE:HD12	1.86	0.41
1:A:24:LEU:HD13	1:A:155:HIS:HE1	1.85	0.41
1:B:112:LEU:HD23	1:B:113:LEU:HD23	2.02	0.41
1:A:164:PHE:CE2	1:A:170:TYR:HB2	2.56	0.41
1:A:203:ARG:HD3	1:A:203:ARG:HA	1.89	0.41
1:C:11:ILE:HG22	1:C:39:TRP:CE2	2.56	0.41
1:C:245:LEU:CD2	1:D:328:LEU:HD12	2.51	0.41
1:D:51:GLN:HE21	1:D:51:GLN:HB2	1.75	0.41
1:B:20:SER:OG	2:B:504:HOH:O	2.22	0.41
1:C:159:PRO:HG2	1:C:164:PHE:HB2	2.04	0.41
1:C:364:MET:HE2	1:C:369:LYS:HE3	2.03	0.41
1:B:364:MET:HE1	1:B:394:GLU:OE1	2.20	0.40
1:C:264:LEU:HD23	1:C:264:LEU:HA	1.79	0.40
1:B:252:LEU:CD1	1:B:292:ARG:NH1	2.84	0.40
1:B:397:SER:OG	1:B:431:MET:HG3	2.21	0.40
1:C:171:ASP:O	1:C:175:GLU:HG3	2.21	0.40
1:D:159:PRO:HG2	1:D:164:PHE:HB2	2.03	0.40
1:A:115:ASP:OD1	1:A:144:ARG:NE	2.48	0.40
1:B:109:VAL:O	1:B:113:LEU:HG	2.21	0.40
1:D:406:ARG:HB3	2:D:523:HOH:O	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:LYS:NZ	1:D:253:LYS:O[1_655]	2.12	0.08

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	446/455 (98%)	434 (97%)	11 (2%)	1 (0%)	43	57
1	B	441/455 (97%)	426 (97%)	14 (3%)	1 (0%)	43	57
1	C	441/455 (97%)	425 (96%)	15 (3%)	1 (0%)	43	57
1	D	441/455 (97%)	427 (97%)	14 (3%)	0	100	100
All	All	1769/1820 (97%)	1712 (97%)	54 (3%)	3 (0%)	43	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	208	SER
1	C	13	PRO
1	A	13	PRO

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	380/383 (99%)	379 (100%)	1 (0%)	86	93
1	B	378/383 (99%)	377 (100%)	1 (0%)	86	93
1	C	378/383 (99%)	377 (100%)	1 (0%)	86	93
1	D	378/383 (99%)	376 (100%)	2 (0%)	81	90
All	All	1514/1532 (99%)	1509 (100%)	5 (0%)	86	93

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	372	VAL
1	B	43	SER
1	C	372	VAL
1	D	253	LYS
1	D	372	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	236	GLN
1	A	257	ASN
1	A	316	ASN
1	A	387	GLN
1	B	51	GLN
1	B	71	GLN
1	B	165	ASN
1	B	236	GLN
1	B	247	GLN
1	B	305	GLN
1	B	337	ASN
1	B	338	GLN
1	C	67	ASN
1	C	189	ASN
1	C	236	GLN
1	C	257	ASN
1	C	308	GLN
1	C	312	HIS
1	C	316	ASN
1	C	327	GLN
1	C	337	ASN
1	C	453	GLN
1	D	312	HIS
1	D	316	ASN
1	D	327	GLN
1	D	337	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

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	450/455 (98%)	-0.22	9 (2%) 65 61	33, 46, 91, 134	0
1	B	447/455 (98%)	-0.06	12 (2%) 56 52	36, 52, 107, 134	0
1	C	447/455 (98%)	-0.06	6 (1%) 75 72	37, 51, 95, 132	0
1	D	447/455 (98%)	-0.11	12 (2%) 56 52	35, 50, 90, 130	0
All	All	1791/1820 (98%)	-0.11	39 (2%) 62 59	33, 50, 98, 134	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	14	PRO	4.9
1	B	245	LEU	3.5
1	C	245	LEU	3.3
1	D	252	LEU	2.9
1	D	351	TYR	2.8
1	C	252	LEU	2.8
1	D	243	PRO	2.7
1	A	252	LEU	2.7
1	B	74	ASP	2.6
1	A	14	PRO	2.6
1	A	245	LEU	2.6
1	C	14	PRO	2.5
1	C	68	LEU	2.5
1	A	21	ALA	2.5
1	A	351	TYR	2.5
1	C	248	LEU	2.5
1	D	209	ALA	2.5
1	B	46	THR	2.5
1	B	299	THR	2.4
1	B	306	ALA	2.3
1	D	245	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	241	LEU	2.2
1	A	304	VAL	2.2
1	D	21	ALA	2.2
1	D	13	PRO	2.2
1	C	9	ASN	2.2
1	D	246	ALA	2.1
1	B	42	TRP	2.1
1	B	304	VAL	2.1
1	B	351	TYR	2.1
1	A	299	THR	2.1
1	A	241	LEU	2.1
1	B	44	GLY	2.1
1	B	300	SER	2.1
1	D	306	ALA	2.0
1	D	53	LEU	2.0
1	A	240	PRO	2.0
1	B	307	TYR	2.0
1	D	456	PRO	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.