



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:17 PM GMT

PDB ID : 1A2O
Title : STRUCTURAL BASIS FOR METHYLESTERASE CHEB REGULATION
BY A PHOSPHORYLATION-ACTIVATED DOMAIN
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Deposited on : 1998-01-06
Resolution : 2.40 Å(reported)

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

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

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

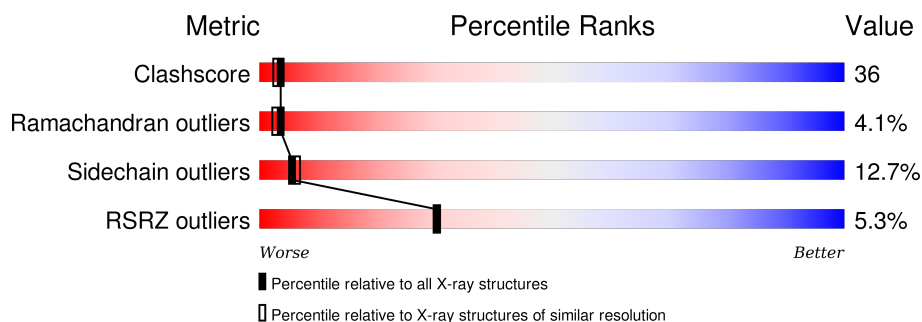
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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

Mol	Chain	Length	Quality of chain
1	A	349	
1	B	349	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5615 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 CHEB METHYLESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	0	0
			2602	1633	465	481	23			
1	B	347	Total	C	N	O	S	0	0	0
			2602	1633	465	481	23			

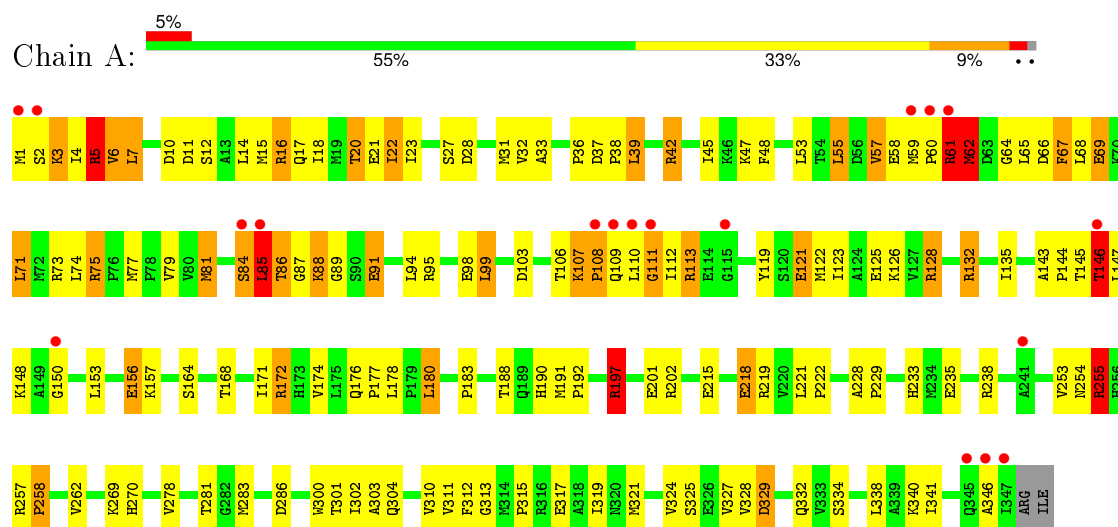
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	221	Total	O	0	0
			221	221		
2	B	190	Total	O	0	0
			190	190		

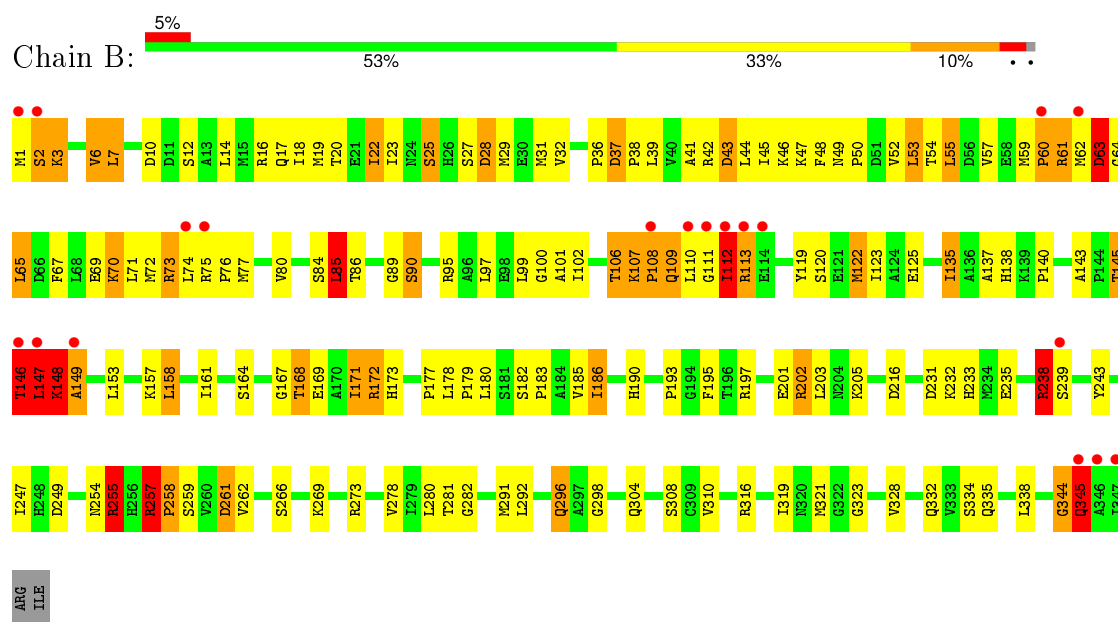
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). 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: CHEB METHYLESTERASE



• Molecule 1: CHEB METHYLESTERASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	163.83 Å 100.46 Å 53.12 Å 90.00° 98.63° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40 14.98 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.9 (10.00-2.40) 97.4 (14.98-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.93 (at 2.20 Å)	Xtriage
Refinement program	REFMAC, X-PLOR	Depositor
R, R_{free}	0.222 , (Not available) 0.229 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.495	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 90.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 41964 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5615	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0.53	0/2645	1.36	31/3577 (0.9%)
1	B	0.52	0/2645	1.30	18/3577 (0.5%)
All	All	0.52	0/5290	1.33	49/7154 (0.7%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	128	ARG	NE-CZ-NH1	15.51	128.05	120.30
1	A	172	ARG	NE-CZ-NH2	-12.73	113.94	120.30
1	A	219	ARG	NE-CZ-NH1	10.88	125.74	120.30
1	B	145	THR	C-N-CA	10.64	148.31	121.70
1	B	172	ARG	NE-CZ-NH2	10.01	125.30	120.30
1	A	172	ARG	NE-CZ-NH1	9.44	125.02	120.30
1	B	257	ARG	NE-CZ-NH2	-8.88	115.86	120.30
1	B	261	ASP	CB-CG-OD1	8.58	126.02	118.30
1	A	128	ARG	CD-NE-CZ	8.45	135.43	123.60
1	B	238	ARG	NE-CZ-NH2	-8.29	116.16	120.30
1	B	243	TYR	CB-CG-CD2	-7.65	116.41	121.00
1	A	197	ARG	NE-CZ-NH1	7.62	124.11	120.30
1	A	197	ARG	NE-CZ-NH2	-7.50	116.55	120.30
1	A	255	ARG	NE-CZ-NH2	-7.38	116.61	120.30
1	A	219	ARG	NE-CZ-NH2	-7.33	116.64	120.30
1	A	257	ARG	CA-C-O	-7.23	104.91	120.10
1	B	172	ARG	NE-CZ-NH1	-7.21	116.69	120.30
1	A	128	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	A	257	ARG	NE-CZ-NH1	6.94	123.77	120.30
1	B	238	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	A	67	PHE	CB-CG-CD1	6.78	125.54	120.80
1	A	42	ARG	NE-CZ-NH1	6.67	123.64	120.30
1	A	172	ARG	CD-NE-CZ	6.63	132.89	123.60
1	B	43	ASP	CB-CG-OD2	6.59	124.23	118.30
1	B	113	ARG	NE-CZ-NH1	-6.39	117.11	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	ARG	CD-NE-CZ	6.33	132.47	123.60
1	A	42	ARG	NE-CZ-NH2	-6.31	117.14	120.30
1	A	329	ASP	CB-CG-OD1	-6.31	112.62	118.30
1	A	255	ARG	NH1-CZ-NH2	5.85	125.83	119.40
1	B	146	THR	N-CA-C	5.85	126.79	111.00
1	B	122	MET	CA-CB-CG	5.77	123.11	113.30
1	A	5	ARG	NE-CZ-NH1	5.65	123.13	120.30
1	B	148	LYS	C-N-CA	5.65	135.81	121.70
1	A	16	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	A	5	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	A	255	ARG	NE-CZ-NH1	-5.48	117.56	120.30
1	B	147	LEU	N-CA-CB	-5.48	99.45	110.40
1	A	91	GLU	OE1-CD-OE2	-5.44	116.77	123.30
1	A	111	GLY	N-CA-C	-5.40	99.61	113.10
1	B	37	ASP	CB-CG-OD1	5.40	123.16	118.30
1	A	67	PHE	CA-CB-CG	5.38	126.81	113.90
1	B	255	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	A	218	GLU	OE1-CD-OE2	-5.37	116.86	123.30
1	A	156	GLU	OE1-CD-OE2	-5.27	116.97	123.30
1	A	67	PHE	N-CA-CB	5.24	120.03	110.60
1	B	63	ASP	CB-CG-OD1	5.18	122.96	118.30
1	B	112	ILE	N-CA-C	5.08	124.70	111.00
1	A	180	LEU	CA-CB-CG	5.04	126.88	115.30
1	A	258	PRO	N-CD-CG	5.02	110.74	103.20

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	2602	0	2686	183	0
1	B	2602	0	2685	196	0
2	A	221	0	0	20	0
2	B	190	0	0	11	0
All	All	5615	0	5371	379	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 36.

All (379) 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:A:319:ILE:HG13	2:A:547:HOH:O	1.40	1.20
1:A:85:LEU:HD11	1:A:108:PRO:HD3	1.21	1.19
1:A:85:LEU:N	1:A:85:LEU:HD22	1.55	1.17
1:B:147:LEU:HD22	1:B:338:LEU:HD23	1.26	1.16
1:B:157:LYS:HZ1	1:B:345:GLN:HG2	1.09	1.14
1:B:109:GLN:NE2	1:B:112:ILE:HG13	1.66	1.11
1:B:147:LEU:HD11	1:B:178:LEU:HD23	1.32	1.07
1:B:147:LEU:HD22	1:B:338:LEU:CD2	1.85	1.06
1:B:110:LEU:O	1:B:110:LEU:HD12	1.55	1.06
1:B:42:ARG:HD3	2:B:506:HOH:O	1.56	1.05
1:A:144:PRO:HB2	1:A:146:THR:OG1	1.56	1.05
1:A:109:GLN:HE22	1:A:110:LEU:HB2	1.24	1.01
1:A:22:ILE:HG13	1:A:23:ILE:N	1.74	1.00
1:A:147:LEU:HD12	1:A:338:LEU:HD22	1.40	1.00
1:B:157:LYS:NZ	1:B:345:GLN:HG2	1.77	0.99
1:A:88:LYS:HA	1:A:88:LYS:HE3	1.41	0.99
1:B:147:LEU:CD2	1:B:338:LEU:CD2	2.40	0.98
1:B:147:LEU:HD21	1:B:338:LEU:HB2	1.46	0.97
1:B:109:GLN:HE21	1:B:112:ILE:HG13	1.26	0.97
1:B:147:LEU:CD1	1:B:178:LEU:HA	1.95	0.96
1:B:147:LEU:HD13	1:B:147:LEU:O	1.67	0.95
1:A:109:GLN:NE2	1:A:110:LEU:HB2	1.80	0.95
1:A:10:ASP:OD2	1:A:15:MET:HG2	1.66	0.94
1:B:149:ALA:HA	2:B:509:HOH:O	1.67	0.94
1:B:110:LEU:O	1:B:110:LEU:CD1	2.15	0.94
1:B:65:LEU:HD12	1:B:99:LEU:HD23	1.51	0.92
1:B:257:ARG:HB3	1:B:258:PRO:HD3	1.50	0.92
1:A:147:LEU:HD11	1:A:178:LEU:HD23	1.53	0.91
1:B:167:GLY:O	1:B:171:ILE:HG23	1.70	0.91
1:B:108:PRO:HG3	1:B:122:MET:CE	2.00	0.91
1:B:147:LEU:O	1:B:177:PRO:O	1.90	0.90
1:A:94:LEU:O	1:A:98:GLU:HG2	1.71	0.89
1:A:5:ARG:HB3	1:A:32:VAL:HG21	1.56	0.88
1:A:22:ILE:HD11	1:A:123:ILE:HG21	1.55	0.87
1:B:119:TYR:CE1	1:B:123:ILE:HD11	2.10	0.87
1:B:22:ILE:HD13	1:B:23:ILE:N	1.89	0.87
1:A:4:ILE:HD11	1:A:132:ARG:NH2	1.90	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:LEU:HD22	1:A:85:LEU:H	1.42	0.85
1:A:300:TRP:CD1	1:A:340:LYS:HG3	2.11	0.85
1:A:65:LEU:O	1:A:69:GLU:HB2	1.75	0.85
1:A:7:LEU:HB3	2:A:535:HOH:O	1.76	0.85
1:B:147:LEU:HD21	1:B:338:LEU:CB	2.07	0.84
1:B:111:GLY:C	1:B:112:ILE:HG12	1.97	0.83
1:A:147:LEU:HD21	1:A:334:SER:OG	1.79	0.82
1:A:238:ARG:HD3	2:A:549:HOH:O	1.80	0.82
1:B:147:LEU:CD1	1:B:178:LEU:HD23	2.08	0.81
1:B:112:ILE:HG22	1:B:113:ARG:H	1.45	0.81
1:A:107:LYS:HD2	1:A:107:LYS:O	1.79	0.81
1:B:147:LEU:CD2	1:B:338:LEU:HD22	2.10	0.81
1:A:85:LEU:HD21	1:A:108:PRO:HA	1.60	0.80
1:A:85:LEU:N	1:A:85:LEU:CD2	2.32	0.80
1:B:145:THR:O	1:B:148:LYS:CE	2.29	0.80
1:B:147:LEU:CD2	1:B:338:LEU:HD23	2.07	0.80
1:A:108:PRO:O	1:A:109:GLN:HB3	1.82	0.80
1:B:55:LEU:HD13	1:B:57:VAL:CG2	2.11	0.80
1:A:147:LEU:HD12	1:A:338:LEU:CD2	2.12	0.79
1:B:158:LEU:HD11	1:B:186:ILE:HD13	1.63	0.78
1:A:55:LEU:HD12	1:A:81:MET:HE1	1.65	0.78
1:B:106:THR:O	1:B:108:PRO:HD3	1.84	0.77
1:B:61:ARG:NE	1:B:61:ARG:HA	2.00	0.77
1:B:147:LEU:CD2	1:B:338:LEU:CB	2.62	0.76
1:B:158:LEU:HD11	1:B:186:ILE:CD1	2.16	0.76
1:B:147:LEU:HD11	1:B:178:LEU:HA	1.67	0.76
1:A:22:ILE:HD11	1:A:123:ILE:CG2	2.16	0.75
1:A:84:SER:C	1:A:85:LEU:HD22	2.05	0.75
1:A:192:PRO:HG3	2:A:417:HOH:O	1.86	0.75
1:A:53:LEU:HD12	1:A:79:VAL:HG13	1.68	0.74
1:B:147:LEU:O	1:B:147:LEU:CD1	2.35	0.74
1:A:85:LEU:HD21	1:A:108:PRO:CA	2.16	0.74
1:A:91:GLU:HB2	2:A:502:HOH:O	1.86	0.74
1:A:17:GLN:OE1	2:A:437:HOH:O	2.06	0.73
1:B:171:ILE:HD11	1:B:203:LEU:CD1	2.18	0.73
1:B:153:LEU:HD12	1:B:183:PRO:HG3	1.70	0.72
1:B:69:GLU:O	1:B:73:ARG:HB2	1.89	0.72
1:A:61:ARG:O	1:A:62:MET:C	2.28	0.72
1:B:147:LEU:HD21	1:B:338:LEU:CD2	2.17	0.72
1:B:257:ARG:HB3	1:B:258:PRO:CD	2.19	0.72
1:B:108:PRO:HG3	1:B:122:MET:HE1	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:LYS:HA	1:A:88:LYS:CE	2.16	0.72
1:B:135:ILE:N	1:B:135:ILE:HD13	2.04	0.72
1:A:37:ASP:HB2	1:A:38:PRO:HD2	1.70	0.72
1:B:147:LEU:HD21	1:B:338:LEU:HD22	1.70	0.71
1:B:7:LEU:O	1:B:53:LEU:HD22	1.90	0.71
1:A:85:LEU:CD1	1:A:108:PRO:HD3	2.11	0.71
1:A:221:LEU:HB3	1:A:222:PRO:HD2	1.73	0.71
1:A:69:GLU:O	1:A:73:ARG:HG3	1.89	0.70
1:A:88:LYS:HE2	1:A:91:GLU:HB3	1.73	0.70
1:B:85:LEU:H	1:B:85:LEU:HD22	1.57	0.70
1:B:147:LEU:HD11	1:B:178:LEU:CD2	2.16	0.70
1:A:147:LEU:H	1:A:147:LEU:HD23	1.56	0.70
1:A:5:ARG:HB3	1:A:32:VAL:CG2	2.22	0.69
1:A:62:MET:HG2	2:A:512:HOH:O	1.92	0.69
1:B:95:ARG:HD3	1:B:202:ARG:HH22	1.58	0.69
1:B:158:LEU:CD1	1:B:186:ILE:HD13	2.22	0.69
1:B:20:THR:HG22	1:B:31:MET:HG2	1.74	0.69
1:A:45:ILE:HG21	1:A:77:MET:CE	2.22	0.69
1:A:11:ASP:HB2	1:A:59:MET:HG2	1.75	0.68
1:A:45:ILE:CG2	1:A:77:MET:HE1	2.24	0.68
1:A:55:LEU:HD12	1:A:81:MET:CE	2.24	0.67
1:B:235:GLU:HB2	1:B:266:SER:OG	1.92	0.67
1:B:257:ARG:O	1:B:259:SER:N	2.28	0.67
1:B:41:ALA:O	1:B:45:ILE:HG13	1.95	0.67
1:B:292:LEU:O	1:B:296:GLN:HG2	1.95	0.67
1:A:147:LEU:CD1	1:A:178:LEU:HD23	2.25	0.67
1:B:110:LEU:C	1:B:110:LEU:HD12	2.15	0.66
1:A:310:VAL:HG22	1:A:311:VAL:HG23	1.77	0.66
1:B:161:ILE:HG12	1:B:278:VAL:HB	1.76	0.66
1:B:32:VAL:HG22	1:B:48:PHE:CD2	2.31	0.66
1:A:281:THR:H	1:A:304:GLN:NE2	1.93	0.66
1:A:147:LEU:CD2	1:A:177:PRO:HB2	2.26	0.66
1:B:153:LEU:HD22	1:B:345:GLN:HG3	1.78	0.65
1:B:22:ILE:HD13	1:B:23:ILE:H	1.56	0.65
1:A:10:ASP:OD2	1:A:15:MET:CG	2.43	0.64
1:B:147:LEU:HD13	1:B:147:LEU:C	2.17	0.64
1:B:62:MET:HB3	2:B:507:HOH:O	1.97	0.64
1:A:110:LEU:C	1:A:112:ILE:H	2.00	0.63
1:A:39:LEU:CD1	1:A:42:ARG:HH21	2.12	0.63
1:A:110:LEU:HG	1:A:111:GLY:H	1.64	0.62
1:B:7:LEU:HD21	1:B:44:LEU:HB2	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:LEU:HD23	1:B:335:GLN:HA	1.79	0.62
1:B:135:ILE:CD1	1:B:135:ILE:H	2.12	0.62
1:B:108:PRO:O	1:B:109:GLN:O	2.18	0.62
1:B:147:LEU:CD2	1:B:338:LEU:HB2	2.22	0.62
1:A:128:ARG:O	1:A:132:ARG:HD2	2.00	0.62
1:B:171:ILE:HD11	1:B:203:LEU:HD11	1.81	0.61
1:A:77:MET:HE2	2:A:524:HOH:O	2.01	0.61
1:B:171:ILE:HD11	1:B:203:LEU:HD13	1.83	0.61
1:B:16:ARG:O	1:B:20:THR:HG23	2.01	0.61
1:B:73:ARG:HH22	1:B:140:PRO:HG2	1.67	0.60
1:B:135:ILE:CD1	1:B:135:ILE:N	2.64	0.60
1:A:147:LEU:C	1:A:148:LYS:HG2	2.22	0.60
1:A:2:SER:HA	1:A:28:ASP:C	2.21	0.60
1:A:321:MET:HG3	2:A:369:HOH:O	2.01	0.60
1:A:147:LEU:HD21	1:A:177:PRO:HB2	1.82	0.60
1:B:61:ARG:NH1	1:B:62:MET:HG2	2.16	0.60
1:B:147:LEU:HD12	1:B:178:LEU:HA	1.83	0.60
1:A:53:LEU:CD1	1:A:79:VAL:HG13	2.32	0.60
1:B:216:ASP:HB2	1:B:247:ILE:HB	1.84	0.60
1:A:45:ILE:HG21	1:A:77:MET:HE1	1.82	0.60
1:A:301:THR:OG1	1:A:324:VAL:HA	2.02	0.59
1:B:112:ILE:CG2	1:B:113:ARG:H	2.07	0.59
1:A:3:LYS:NZ	1:A:28:ASP:HB2	2.17	0.59
1:A:39:LEU:HD13	1:A:42:ARG:HH21	1.67	0.59
1:B:205:LYS:HD3	2:B:459:HOH:O	2.01	0.59
1:A:328:VAL:HG12	1:A:329:ASP:O	2.03	0.59
1:A:283:MET:HE2	2:A:553:HOH:O	2.01	0.59
1:A:55:LEU:CD1	1:A:81:MET:CE	2.81	0.59
1:B:119:TYR:CZ	1:B:123:ILE:HD11	2.38	0.59
1:A:55:LEU:CD1	1:A:81:MET:HE3	2.33	0.58
1:A:88:LYS:CE	1:A:91:GLU:HB3	2.33	0.58
1:B:145:THR:O	1:B:148:LYS:HE3	2.03	0.58
1:A:95:ARG:HD2	1:A:202:ARG:HH22	1.68	0.58
1:B:61:ARG:HA	1:B:61:ARG:HE	1.67	0.58
1:A:85:LEU:HD13	1:A:85:LEU:H	1.67	0.58
1:B:57:VAL:HG13	1:B:64:GLY:HA3	1.85	0.57
1:A:183:PRO:HG2	1:A:341:ILE:HG21	1.87	0.57
1:B:147:LEU:HD12	1:B:177:PRO:C	2.24	0.57
1:B:233:HIS:CE1	1:B:257:ARG:O	2.58	0.57
1:A:37:ASP:OD1	1:A:39:LEU:HB2	2.05	0.57
1:A:10:ASP:CG	1:A:15:MET:HG2	2.25	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ASP:HB2	1:B:38:PRO:CD	2.35	0.57
1:A:84:SER:O	1:A:85:LEU:C	2.43	0.56
1:B:157:LYS:HZ1	1:B:345:GLN:HA	1.69	0.56
1:A:45:ILE:CG2	1:A:77:MET:CE	2.83	0.56
1:A:319:ILE:CG1	2:A:547:HOH:O	2.20	0.56
1:A:147:LEU:O	1:A:148:LYS:HG2	2.05	0.56
1:A:303:ALA:HB3	1:A:319:ILE:HD11	1.87	0.56
1:B:122:MET:O	1:B:125:GLU:HG3	2.06	0.56
1:A:22:ILE:HG13	1:A:23:ILE:H	1.70	0.56
1:A:176:GLN:HB3	1:A:177:PRO:HD3	1.87	0.56
1:B:157:LYS:HZ1	1:B:345:GLN:CG	1.99	0.55
1:A:55:LEU:HB3	1:A:81:MET:HE3	1.89	0.55
1:A:74:LEU:O	1:A:75:ARG:HB2	2.05	0.55
1:A:328:VAL:CG1	1:A:332:GLN:HB2	2.36	0.55
1:A:32:VAL:O	1:A:33:ALA:HB2	2.06	0.55
1:B:100:GLY:O	1:B:138:HIS:NE2	2.39	0.54
1:B:147:LEU:C	1:B:148:LYS:CG	2.76	0.54
1:B:97:LEU:HA	1:B:101:ALA:HB3	1.90	0.54
1:B:10:ASP:O	1:B:36:PRO:HA	2.07	0.54
1:A:192:PRO:CG	2:A:417:HOH:O	2.51	0.54
1:A:95:ARG:NH2	2:A:475:HOH:O	2.39	0.54
1:A:312:PHE:O	1:A:315:PRO:HD2	2.08	0.54
1:B:106:THR:O	1:B:108:PRO:CD	2.56	0.54
1:B:45:ILE:HD13	1:B:77:MET:HE3	1.90	0.54
1:B:119:TYR:O	1:B:123:ILE:HD12	2.08	0.54
1:A:71:LEU:HD22	1:A:71:LEU:O	2.08	0.54
1:B:147:LEU:HG	1:B:334:SER:OG	2.08	0.54
1:B:108:PRO:HG3	1:B:122:MET:HE3	1.88	0.54
1:A:55:LEU:HD13	1:A:81:MET:HE3	1.88	0.54
1:B:2:SER:HB3	2:B:424:HOH:O	2.06	0.54
1:B:145:THR:O	1:B:148:LYS:CD	2.55	0.54
1:A:190:HIS:NE2	1:A:286:ASP:OD2	2.41	0.54
1:A:150:GLY:N	2:A:364:HOH:O	2.41	0.53
1:A:3:LYS:HE2	1:A:132:ARG:NH2	2.24	0.53
1:A:147:LEU:HD11	1:A:178:LEU:CD2	2.34	0.53
1:B:89:GLY:O	1:B:90:SER:CB	2.55	0.53
1:B:231:ASP:C	1:B:257:ARG:HG2	2.29	0.53
1:A:346:ALA:HB2	2:A:488:HOH:O	2.09	0.53
1:B:112:ILE:HG22	1:B:113:ARG:N	2.20	0.53
1:A:16:ARG:O	1:A:20:THR:HG23	2.08	0.53
1:B:147:LEU:CD2	1:B:338:LEU:HB3	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:MET:O	1:B:63:ASP:O	2.27	0.53
1:A:110:LEU:C	1:A:112:ILE:N	2.62	0.53
1:B:145:THR:O	1:B:148:LYS:HE2	2.08	0.53
1:A:64:GLY:O	1:A:68:LEU:HB3	2.09	0.52
1:A:145:THR:OG1	1:A:146:THR:N	2.42	0.52
1:A:4:ILE:HD11	1:A:132:ARG:HH21	1.71	0.52
1:B:61:ARG:O	1:B:62:MET:HB2	2.08	0.52
1:B:254:ASN:OD1	1:B:261:ASP:OD2	2.25	0.52
1:B:157:LYS:NZ	1:B:345:GLN:HA	2.25	0.52
1:A:58:GLU:HG2	1:A:58:GLU:O	2.10	0.52
1:B:32:VAL:HG22	1:B:48:PHE:HD2	1.72	0.52
1:A:147:LEU:O	1:A:177:PRO:O	2.28	0.52
1:A:3:LYS:HE2	1:A:132:ARG:HH21	1.75	0.51
1:B:147:LEU:O	1:B:148:LYS:HG2	2.09	0.51
1:B:44:LEU:N	1:B:44:LEU:HD23	2.25	0.51
1:B:95:ARG:O	1:B:99:LEU:HB2	2.11	0.51
1:A:113:ARG:HA	1:A:113:ARG:HE	1.75	0.51
1:B:84:SER:O	1:B:86:THR:N	2.44	0.51
1:A:84:SER:O	1:A:86:THR:N	2.43	0.51
1:B:193:PRO:HG3	1:B:231:ASP:HB2	1.93	0.51
1:A:119:TYR:CZ	1:A:123:ILE:HD11	2.45	0.51
1:B:232:LYS:HD3	1:B:249:ASP:OD2	2.10	0.51
1:A:47:LYS:HG2	1:A:48:PHE:CD2	2.46	0.51
1:B:19:MET:O	1:B:22:ILE:HD13	2.11	0.51
1:A:215:GLU:HG2	1:A:218:GLU:OE2	2.10	0.51
1:B:238:ARG:NH2	2:B:378:HOH:O	2.44	0.50
1:A:32:VAL:HG11	1:A:48:PHE:HB3	1.92	0.50
1:B:18:ILE:HD13	1:B:18:ILE:N	2.27	0.50
1:B:10:ASP:HB3	1:B:16:ARG:HG3	1.92	0.50
1:A:269:LYS:HE2	1:A:270:HIS:CE1	2.47	0.50
1:B:65:LEU:CD1	1:B:99:LEU:HD23	2.32	0.50
1:A:6:VAL:HG22	1:A:31:MET:SD	2.51	0.50
1:B:17:GLN:O	1:B:20:THR:OG1	2.28	0.50
1:B:2:SER:HA	1:B:29:MET:N	2.26	0.50
1:A:109:GLN:OE1	1:A:112:ILE:HG21	2.11	0.50
1:B:62:MET:O	1:B:63:ASP:C	2.49	0.50
1:A:255:ARG:HA	1:A:255:ARG:NE	2.26	0.49
1:A:171:ILE:O	1:A:174:VAL:HG22	2.12	0.49
1:A:18:ILE:O	1:A:22:ILE:HG23	2.13	0.49
1:A:321:MET:CG	2:A:369:HOH:O	2.60	0.49
1:A:53:LEU:HD12	1:A:79:VAL:CG1	2.40	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:ARG:HD2	1:A:202:ARG:NH2	2.26	0.49
1:B:119:TYR:CD1	1:B:123:ILE:HD11	2.48	0.49
1:A:2:SER:HA	1:A:28:ASP:CA	2.43	0.48
1:A:64:GLY:O	1:A:68:LEU:CB	2.61	0.48
1:B:145:THR:OG1	1:B:146:THR:N	2.45	0.48
1:B:42:ARG:HG3	1:B:43:ASP:N	2.28	0.48
1:A:110:LEU:HG	1:A:111:GLY:N	2.29	0.48
1:B:233:HIS:ND1	1:B:262:VAL:HG21	2.28	0.48
1:A:319:ILE:HD13	1:A:324:VAL:HG11	1.96	0.48
1:B:147:LEU:HD13	1:B:179:PRO:HD3	1.95	0.48
1:A:84:SER:CA	1:A:85:LEU:HD22	2.44	0.47
1:B:110:LEU:O	1:B:110:LEU:HD13	2.11	0.47
1:B:135:ILE:H	1:B:135:ILE:HD13	1.70	0.47
1:A:57:VAL:HG23	2:A:501:HOH:O	2.13	0.47
1:B:55:LEU:HD13	1:B:57:VAL:HG22	1.93	0.47
1:A:153:LEU:HD12	1:A:183:PRO:HG3	1.96	0.47
1:B:135:ILE:HA	1:B:138:HIS:HB3	1.96	0.47
1:A:188:THR:HG23	1:A:258:PRO:HB2	1.96	0.47
1:A:109:GLN:CD	1:A:110:LEU:HB2	2.35	0.47
1:A:16:ARG:NH1	1:A:36:PRO:HB3	2.29	0.47
1:B:74:LEU:C	1:B:76:PRO:HD3	2.33	0.47
1:A:233:HIS:CE1	1:A:262:VAL:HG21	2.49	0.47
1:A:32:VAL:CG1	1:A:48:PHE:HB3	2.44	0.47
1:B:344:GLY:O	1:B:345:GLN:HB2	2.14	0.47
1:A:281:THR:H	1:A:304:GLN:HE22	1.63	0.47
1:A:253:VAL:HG12	1:A:254:ASN:OD1	2.15	0.46
1:B:95:ARG:CD	1:B:202:ARG:HH22	2.28	0.46
1:B:61:ARG:CA	1:B:61:ARG:NE	2.75	0.46
1:A:310:VAL:CG2	1:A:311:VAL:HG23	2.43	0.46
1:A:16:ARG:CZ	1:A:36:PRO:HB3	2.45	0.46
1:A:319:ILE:HD12	1:A:327:VAL:HG22	1.98	0.46
1:B:42:ARG:HG3	1:B:43:ASP:OD1	2.16	0.46
1:B:45:ILE:HD13	1:B:77:MET:CE	2.44	0.46
1:A:2:SER:HA	1:A:28:ASP:HA	1.97	0.46
1:A:215:GLU:O	1:A:218:GLU:HG2	2.16	0.46
1:A:319:ILE:CD1	1:A:324:VAL:HG11	2.46	0.46
1:B:107:LYS:O	1:B:108:PRO:C	2.54	0.46
1:A:16:ARG:NH2	1:A:36:PRO:HB3	2.31	0.46
1:B:255:ARG:HA	1:B:255:ARG:HD3	1.80	0.46
1:A:16:ARG:O	1:A:20:THR:CG2	2.64	0.46
1:A:45:ILE:HG23	1:A:77:MET:HE1	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:LYS:HD3	2:A:545:HOH:O	2.16	0.46
1:B:255:ARG:CD	1:B:255:ARG:N	2.79	0.46
1:A:103:ASP:OD2	1:A:126:LYS:NZ	2.48	0.46
1:B:72:MET:HG3	1:B:100:GLY:HA3	1.98	0.46
1:A:107:LYS:O	1:A:107:LYS:CD	2.59	0.45
1:B:32:VAL:CG2	1:B:48:PHE:HD2	2.28	0.45
1:A:85:LEU:CD2	1:A:85:LEU:H	2.10	0.45
1:B:38:PRO:HB3	1:B:67:PHE:HB2	1.98	0.45
1:B:195:PHE:HD2	2:B:521:HOH:O	1.99	0.45
1:A:228:ALA:HA	1:A:229:PRO:HD3	1.75	0.45
1:B:197:ARG:NH1	1:B:201:GLU:OE1	2.49	0.45
1:B:172:ARG:HG3	1:B:203:LEU:HD23	1.99	0.45
1:B:49:ASN:N	1:B:50:PRO:HD3	2.31	0.45
1:A:317:GLU:O	1:A:321:MET:HG3	2.17	0.45
1:B:90:SER:HA	1:B:195:PHE:CE2	2.51	0.45
1:B:113:ARG:N	1:B:113:ARG:HD2	2.32	0.44
1:B:257:ARG:CB	1:B:258:PRO:CD	2.91	0.44
1:B:60:PRO:HB2	1:B:61:ARG:H	1.61	0.44
1:A:253:VAL:O	1:A:254:ASN:HB2	2.17	0.44
1:B:316:ARG:HA	1:B:319:ILE:HD12	1.99	0.44
1:B:54:THR:HA	1:B:80:VAL:O	2.17	0.44
1:B:257:ARG:O	1:B:258:PRO:C	2.55	0.44
1:B:135:ILE:O	1:B:138:HIS:HB3	2.16	0.44
1:A:135:ILE:N	1:A:135:ILE:HD12	2.33	0.44
1:B:27:SER:O	1:B:28:ASP:HB3	2.17	0.44
1:B:6:VAL:HA	1:B:52:VAL:O	2.17	0.44
1:B:110:LEU:C	1:B:110:LEU:CD1	2.82	0.44
1:A:340:LYS:HD3	1:A:340:LYS:HA	1.78	0.44
1:A:122:MET:HE2	1:A:122:MET:HA	1.98	0.44
1:A:85:LEU:CD1	1:A:106:THR:HB	2.47	0.44
1:B:101:ALA:O	1:B:138:HIS:HE1	2.01	0.44
1:B:232:LYS:N	1:B:257:ARG:HG2	2.33	0.44
1:A:73:ARG:NH1	1:A:73:ARG:HG2	2.33	0.44
1:A:147:LEU:HD23	1:A:177:PRO:HB2	2.00	0.44
1:A:176:GLN:N	1:A:177:PRO:CD	2.81	0.43
1:B:61:ARG:HB2	2:B:423:HOH:O	2.18	0.43
1:A:283:MET:CE	2:A:553:HOH:O	2.64	0.43
1:B:71:LEU:CD2	1:B:75:ARG:O	2.66	0.43
1:B:168:THR:HG22	1:B:169:GLU:N	2.33	0.43
1:B:112:ILE:CG2	1:B:113:ARG:N	2.79	0.43
1:B:19:MET:O	1:B:22:ILE:CD1	2.65	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:TRP:NE1	1:A:340:LYS:HG3	2.33	0.43
1:A:73:ARG:HH11	1:A:73:ARG:HG2	1.83	0.43
1:A:319:ILE:CD1	1:A:327:VAL:HG22	2.48	0.43
1:A:10:ASP:OD2	1:A:15:MET:SD	2.77	0.43
1:B:344:GLY:O	1:B:345:GLN:CB	2.67	0.43
1:A:132:ARG:HD2	1:A:132:ARG:N	2.33	0.43
1:A:221:LEU:CB	1:A:222:PRO:HD2	2.44	0.43
1:A:143:ALA:HB2	1:A:172:ARG:NH2	2.34	0.43
1:A:121:GLU:O	1:A:125:GLU:HG2	2.18	0.43
1:B:108:PRO:HG3	1:B:122:MET:SD	2.59	0.42
1:B:3:LYS:HA	2:B:460:HOH:O	2.19	0.42
1:B:39:LEU:HA	1:B:39:LEU:HD12	1.82	0.42
1:B:273:ARG:HB3	1:B:298:GLY:HA3	2.02	0.42
1:B:179:PRO:HD2	1:B:182:SER:OG	2.19	0.42
1:A:191:MET:HG2	1:A:192:PRO:HD2	2.02	0.42
1:B:37:ASP:HB2	1:B:38:PRO:HD2	2.01	0.42
1:A:238:ARG:HB3	1:A:238:ARG:HE	1.75	0.42
1:A:157:LYS:HD3	1:A:157:LYS:N	2.35	0.42
1:B:258:PRO:HD2	2:B:368:HOH:O	2.19	0.42
1:A:156:GLU:OE2	1:A:238:ARG:NH1	2.52	0.42
1:A:278:VAL:HA	1:A:302:ILE:O	2.18	0.42
1:B:190:HIS:HA	2:B:368:HOH:O	2.20	0.42
1:A:1:MET:HG2	1:A:27:SER:O	2.20	0.42
1:B:269:LYS:HE3	1:B:269:LYS:HB2	1.89	0.41
1:B:147:LEU:HD23	1:B:338:LEU:HB3	2.00	0.41
1:B:73:ARG:HH22	1:B:140:PRO:CG	2.32	0.41
1:A:39:LEU:HA	1:A:39:LEU:HD13	1.77	0.41
1:B:143:ALA:HA	1:B:173:HIS:CD2	2.55	0.41
1:B:280:LEU:HB3	1:B:304:GLN:NE2	2.36	0.41
1:B:292:LEU:CD2	1:B:321:MET:HE2	2.50	0.41
1:B:328:VAL:CG1	1:B:332:GLN:HB3	2.51	0.41
1:A:197:ARG:NH1	1:A:201:GLU:OE1	2.52	0.41
1:A:85:LEU:HD21	1:A:108:PRO:CB	2.50	0.41
1:B:70:LYS:HA	1:B:70:LYS:HD2	1.46	0.41
1:B:147:LEU:HD21	1:B:338:LEU:CG	2.50	0.41
1:A:69:GLU:HG2	1:A:99:LEU:HD21	2.01	0.41
1:A:7:LEU:CD1	1:A:7:LEU:C	2.89	0.41
1:B:55:LEU:HA	1:B:55:LEU:HD23	1.82	0.41
1:B:137:ALA:HB1	1:B:308:SER:O	2.20	0.41
1:A:109:GLN:CD	1:A:109:GLN:C	2.78	0.41
1:B:233:HIS:ND1	1:B:257:ARG:O	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:SER:N	1:A:27:SER:O	2.54	0.40
1:B:100:GLY:O	1:B:138:HIS:CE1	2.75	0.40
1:B:1:MET:O	1:B:2:SER:C	2.57	0.40
1:A:86:THR:O	1:A:89:GLY:N	2.54	0.40
1:A:5:ARG:HG3	2:A:433:HOH:O	2.20	0.40
1:A:312:PHE:O	1:A:313:GLY:C	2.59	0.40
1:B:55:LEU:HD13	1:B:57:VAL:HG21	2.00	0.40
1:B:71:LEU:HD22	1:B:75:ARG:O	2.22	0.40
1:B:164:SER:HB3	1:B:282:GLY:CA	2.51	0.40
1:B:43:ASP:O	1:B:47:LYS:HG3	2.22	0.40
1:B:95:ARG:HD3	1:B:202:ARG:NH2	2.33	0.40
1:B:291:MET:HG2	1:B:323:GLY:O	2.21	0.40
1:B:65:LEU:HA	1:B:65:LEU:HD13	1.73	0.40
1:B:158:LEU:HD11	1:B:186:ILE:HD11	1.97	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	345/349 (99%)	320 (93%)	15 (4%)	10 (3%)	6	5
1	B	345/349 (99%)	300 (87%)	27 (8%)	18 (5%)	2	1
All	All	690/698 (99%)	620 (90%)	42 (6%)	28 (4%)	3	3

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	85	LEU
1	A	108	PRO
1	B	25	SER
1	B	63	ASP

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Mol	Chain	Res	Type
1	B	109	GLN
1	B	112	ILE
1	B	146	THR
1	A	60	PRO
1	A	62	MET
1	A	87	GLY
1	A	146	THR
1	B	28	ASP
1	B	60	PRO
1	B	90	SER
1	B	281	THR
1	B	345	GLN
1	A	86	THR
1	B	61	ARG
1	B	85	LEU
1	B	108	PRO
1	B	149	ALA
1	B	2	SER
1	A	61	ARG
1	B	257	ARG
1	A	164	SER
1	B	258	PRO
1	B	344	GLY
1	A	75	ARG

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	280/282 (99%)	246 (88%)	34 (12%)	6	8
1	B	280/282 (99%)	243 (87%)	37 (13%)	5	6
All	All	560/564 (99%)	489 (87%)	71 (13%)	5	6

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	5	ARG
1	A	6	VAL
1	A	7	LEU
1	A	12	SER
1	A	14	LEU
1	A	20	THR
1	A	21	GLU
1	A	22	ILE
1	A	39	LEU
1	A	55	LEU
1	A	57	VAL
1	A	61	ARG
1	A	62	MET
1	A	66	ASP
1	A	67	PHE
1	A	69	GLU
1	A	71	LEU
1	A	81	MET
1	A	84	SER
1	A	85	LEU
1	A	88	LYS
1	A	99	LEU
1	A	107	LYS
1	A	113	ARG
1	A	121	GLU
1	A	132	ARG
1	A	146	THR
1	A	168	THR
1	A	180	LEU
1	A	197	ARG
1	A	235	GLU
1	A	255	ARG
1	A	325	SER
1	B	3	LYS
1	B	6	VAL
1	B	7	LEU
1	B	12	SER
1	B	14	LEU
1	B	22	ILE
1	B	25	SER
1	B	46	LYS
1	B	53	LEU

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Mol	Chain	Res	Type
1	B	55	LEU
1	B	59	MET
1	B	65	LEU
1	B	70	LYS
1	B	73	ARG
1	B	85	LEU
1	B	102	ILE
1	B	106	THR
1	B	107	LYS
1	B	112	ILE
1	B	120	SER
1	B	135	ILE
1	B	146	THR
1	B	147	LEU
1	B	148	LYS
1	B	158	LEU
1	B	168	THR
1	B	171	ILE
1	B	180	LEU
1	B	185	VAL
1	B	186	ILE
1	B	202	ARG
1	B	238	ARG
1	B	239	SER
1	B	255	ARG
1	B	296	GLN
1	B	310	VAL
1	B	345	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	242	ASN
1	A	304	GLN
1	B	109	GLN
1	B	176	GLN
1	B	254	ASN
1	B	270	HIS
1	B	296	GLN
1	B	304	GLN

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 carbohydrates 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	347/349 (99%)	-0.23	18 (5%) 31 31	6, 20, 54, 60	0
1	B	347/349 (99%)	-0.14	19 (5%) 29 29	6, 23, 50, 60	0
All	All	694/698 (99%)	-0.18	37 (5%) 30 30	6, 22, 52, 60	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	347	ILE	6.2
1	B	111	GLY	6.0
1	B	112	ILE	5.9
1	B	346	ALA	5.9
1	A	110	LEU	4.9
1	A	2	SER	4.9
1	A	111	GLY	4.8
1	A	347	ILE	4.5
1	A	146	THR	4.1
1	B	1	MET	3.9
1	A	1	MET	3.8
1	B	2	SER	3.7
1	A	115	GLY	3.6
1	B	147	LEU	3.4
1	A	61	ARG	3.2
1	B	60	PRO	3.2
1	B	345	GLN	3.2
1	A	84	SER	3.2
1	B	114	GLU	3.1
1	B	62	MET	3.1
1	A	59	MET	3.1
1	B	110	LEU	3.0
1	A	345	GLN	2.9
1	A	85	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	74	LEU	2.8
1	B	113	ARG	2.7
1	A	346	ALA	2.6
1	B	239	SER	2.5
1	A	241	ALA	2.5
1	A	109	GLN	2.5
1	B	146	THR	2.4
1	A	150	GLY	2.4
1	A	60	PRO	2.4
1	A	108	PRO	2.3
1	B	108	PRO	2.3
1	B	149	ALA	2.3
1	B	75	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.