



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:12 AM UTC

PDB ID : 3KJX / pdb_00003kix
Title : Crystal structure of a transcriptional regulator, LacI family protein from *Silicibacter pomeroyi*
Authors : Palani, K.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-11-03
Resolution : 2.33 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

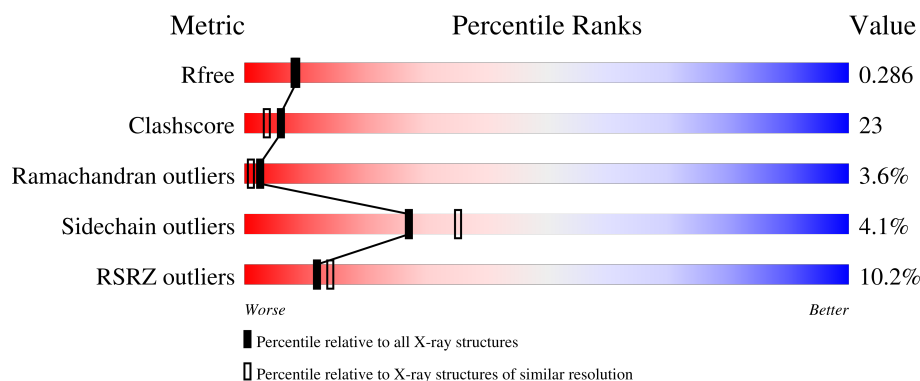
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	344	13% 58% 33% ...
1	B	344	7% 63% 28% 5% ...
1	C	344	8% 53% 37% 7% ...
1	D	344	10% 58% 35% ...

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10352 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 Transcriptional regulator, LacI family.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	Se	0	0	0
			2552	1603	448	487	1	13			
1	B	333	Total	C	N	O	S	Se	0	0	0
			2552	1603	448	487	1	13			
1	C	333	Total	C	N	O	S	Se	0	0	0
			2552	1603	448	487	1	13			
1	D	335	Total	C	N	O	S	Se	0	0	0
			2572	1615	454	489	1	13			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP Q5LM80
A	0	SER	-	expression tag	UNP Q5LM80
A	1	LEU	-	expression tag	UNP Q5LM80
A	341	GLU	-	expression tag	UNP Q5LM80
A	342	GLY	-	expression tag	UNP Q5LM80
B	-1	MSE	-	expression tag	UNP Q5LM80
B	0	SER	-	expression tag	UNP Q5LM80
B	1	LEU	-	expression tag	UNP Q5LM80
B	341	GLU	-	expression tag	UNP Q5LM80
B	342	GLY	-	expression tag	UNP Q5LM80
C	-1	MSE	-	expression tag	UNP Q5LM80
C	0	SER	-	expression tag	UNP Q5LM80
C	1	LEU	-	expression tag	UNP Q5LM80
C	341	GLU	-	expression tag	UNP Q5LM80
C	342	GLY	-	expression tag	UNP Q5LM80
D	-1	MSE	-	expression tag	UNP Q5LM80
D	0	SER	-	expression tag	UNP Q5LM80
D	1	LEU	-	expression tag	UNP Q5LM80
D	341	GLU	-	expression tag	UNP Q5LM80
D	342	GLY	-	expression tag	UNP Q5LM80

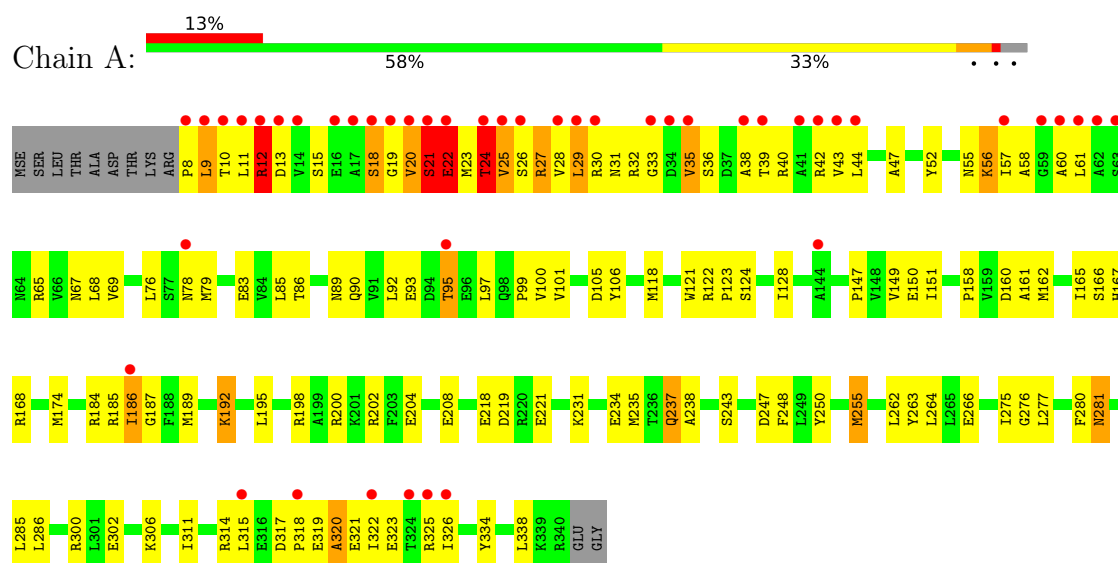
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	36	Total 36	O 36	0	0
2	B	25	Total 25	O 25	0	0
2	C	30	Total 30	O 30	0	0
2	D	33	Total 33	O 33	0	0

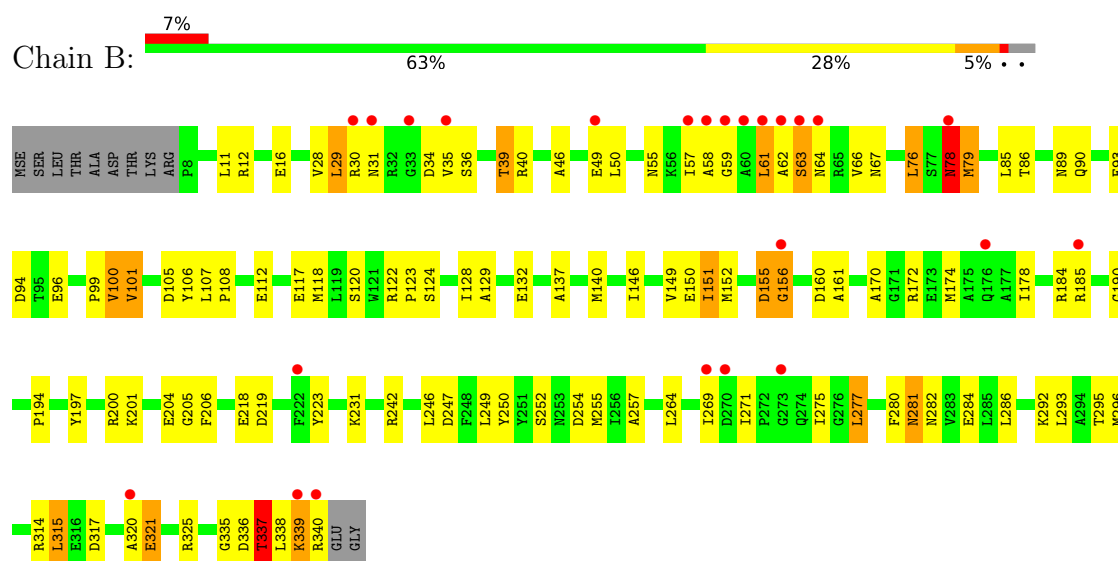
3 Residue-property plots

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.

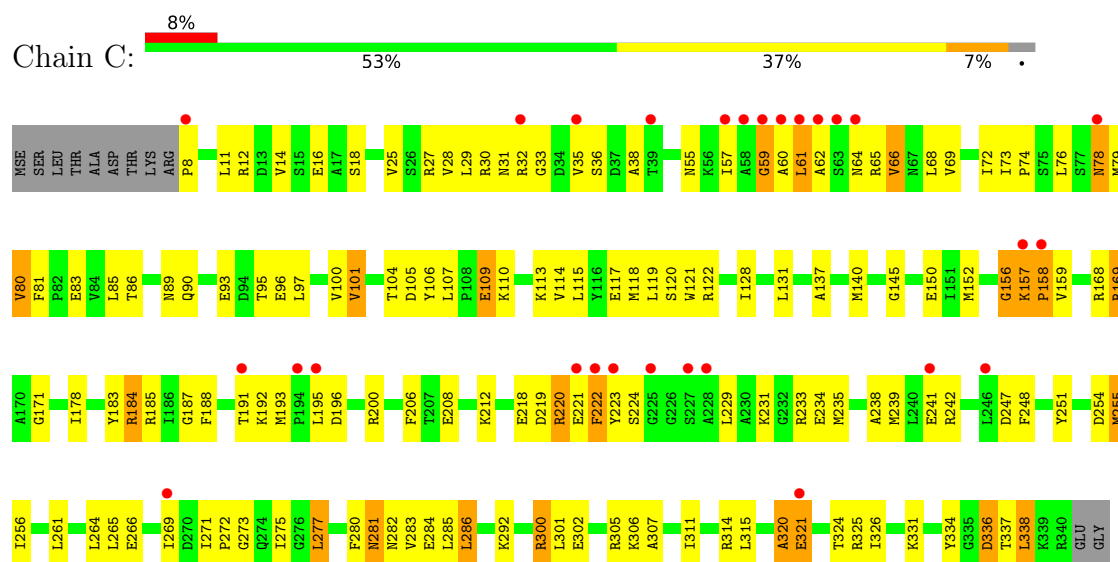
- Molecule 1: Transcriptional regulator, LacI family



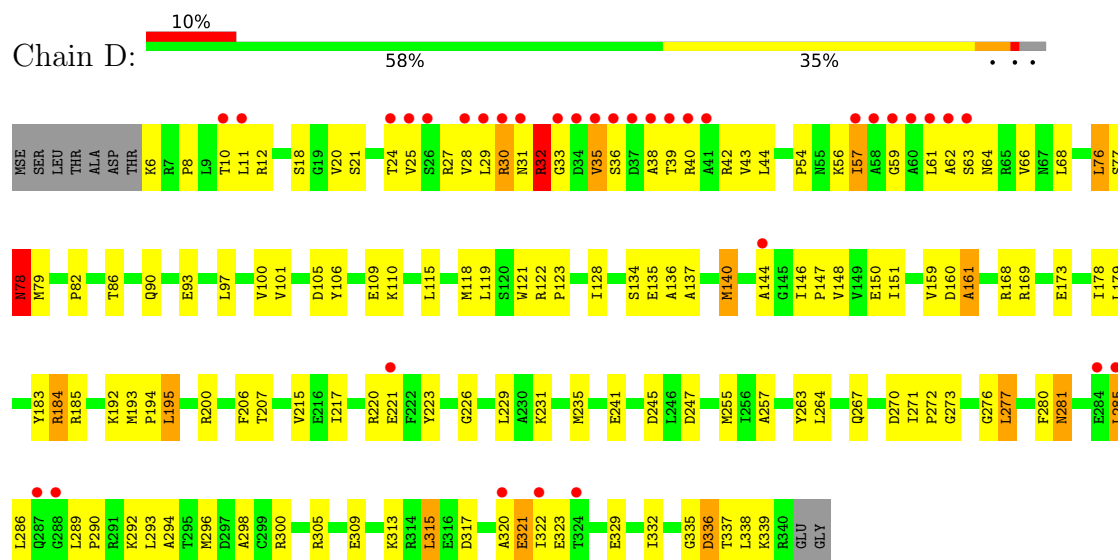
- Molecule 1: Transcriptional regulator, LacI family



- Molecule 1: Transcriptional regulator, LacI family



• Molecule 1: Transcriptional regulator, LacI family



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.66Å 137.39Å 154.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.29 – 2.33 41.29 – 2.33	Depositor EDS
% Data completeness (in resolution range)	94.7 (41.29-2.33) 94.7 (41.29-2.33)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.24Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.238 , 0.287 0.238 , 0.286	Depositor DCC
R_{free} test set	3076 reflections (4.40%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.554	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10352	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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	0.48	1/2574 (0.0%)	0.99	12/3451 (0.3%)
1	B	0.42	0/2574	0.95	6/3451 (0.2%)
1	C	0.46	0/2574	0.97	8/3451 (0.2%)
1	D	0.47	0/2594	1.01	12/3477 (0.3%)
All	All	0.46	1/10316 (0.0%)	0.98	38/13830 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	20	VAL	CA-CB	-8.86	1.43	1.54

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	184	ARG	N-CA-C	-8.68	96.21	109.14
1	D	101	VAL	N-CA-C	8.17	120.04	108.36
1	C	101	VAL	N-CA-C	7.29	119.57	108.71
1	C	78	ASN	N-CA-C	-6.88	100.27	110.46
1	A	243	SER	CA-C-N	6.73	126.42	119.56
1	A	243	SER	C-N-CA	6.73	126.42	119.56
1	B	101	VAL	N-CA-C	6.66	118.19	108.53
1	C	334	TYR	N-CA-C	6.59	120.42	112.38
1	D	151	ILE	N-CA-C	6.40	118.97	108.99
1	D	195	LEU	N-CA-C	6.38	119.05	111.71
1	D	226	GLY	N-CA-C	6.36	120.68	112.54
1	D	159	VAL	N-CA-C	-6.33	104.68	110.82
1	A	10	THR	N-CA-C	6.30	116.59	108.34
1	A	78	ASN	N-CA-C	-6.25	100.43	109.31
1	A	18	SER	N-CA-C	-6.19	104.59	111.71
1	B	76	LEU	N-CA-C	-6.11	104.24	113.89
1	A	334	TYR	N-CA-C	6.09	118.88	111.82
1	B	146	ILE	N-CA-C	6.07	114.58	107.84
1	D	38	ALA	N-CA-C	-6.03	106.53	114.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	151	ILE	N-CA-C	5.97	118.45	109.20
1	A	195	LEU	N-CA-C	5.96	117.85	111.36
1	A	168	ARG	N-CA-C	-5.85	104.81	111.07
1	D	76	LEU	N-CA-C	-5.73	104.83	113.89
1	D	184	ARG	N-CA-C	-5.72	100.62	109.14
1	C	320	ALA	N-CA-C	5.65	117.44	111.28
1	A	151	ILE	N-CA-C	5.65	117.31	109.29
1	D	78	ASN	N-CA-C	-5.64	98.79	110.80
1	B	337	THR	N-CA-C	5.51	122.53	110.80
1	D	136	ALA	N-CA-C	-5.50	105.37	111.36
1	B	34	ASP	N-CA-C	-5.49	104.83	112.30
1	A	21	SER	N-CA-C	5.44	122.39	110.80
1	A	22	GLU	N-CA-C	-5.41	99.28	110.80
1	D	241	GLU	N-CA-C	-5.37	105.51	111.36
1	C	66	VAL	N-CA-C	-5.32	100.19	108.86
1	C	80	VAL	CB-CA-C	-5.16	107.32	111.71
1	D	336	ASP	N-CA-C	-5.10	105.13	113.19
1	C	168	ARG	N-CA-C	-5.04	105.79	111.28
1	A	275	ILE	N-CA-C	5.02	115.35	108.12

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	2552	0	2609	125	0
1	B	2552	0	2609	100	0
1	C	2552	0	2609	143	0
1	D	2572	0	2634	135	0
2	A	36	0	0	2	0
2	B	25	0	0	1	0
2	C	30	0	0	1	0
2	D	33	0	0	3	0
All	All	10352	0	10461	469	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 23.

All (469) 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:C:264:LEU:HG	1:C:269:ILE:HD11	1.24	1.18
1:C:115:LEU:HD23	1:C:140:MSE:HE3	1.40	1.03
1:C:137:ALA:HA	1:C:140:MSE:HE2	1.43	1.00
1:B:249:LEU:HG	1:B:275:ILE:HD11	1.44	0.99
1:B:137:ALA:HA	1:B:140:MSE:HE3	1.45	0.99
1:D:28:VAL:HG21	1:D:43:VAL:HG21	1.45	0.97
1:C:95:THR:HG22	1:C:97:LEU:H	1.27	0.94
1:D:137:ALA:HA	1:D:140:MSE:HE2	1.47	0.94
1:B:277:LEU:HD12	1:B:337:THR:HB	1.52	0.92
1:C:283:VAL:HA	1:C:300:ARG:HH12	1.36	0.90
1:D:20:VAL:HG21	1:D:43:VAL:HG22	1.53	0.89
1:D:294:ALA:H	1:D:337:THR:CG2	1.86	0.88
1:A:255:MSE:HE2	1:A:255:MSE:HA	1.52	0.88
1:D:61:LEU:HA	1:D:122:ARG:HE	1.39	0.86
1:D:322:ILE:HG22	1:D:323:GLU:H	1.39	0.86
1:A:55:ASN:HD22	1:B:61:LEU:H	1.23	0.84
1:A:100:VAL:HB	1:B:100:VAL:HG13	1.60	0.83
1:C:60:ALA:N	1:C:64:ASN:HB2	1.94	0.83
1:C:114:VAL:HG12	1:C:118:MSE:HE2	1.59	0.83
1:D:115:LEU:HD23	1:D:140:MSE:HE3	1.62	0.81
1:A:31:ASN:HA	1:A:40:ARG:HH22	1.44	0.81
1:B:271:ILE:HG22	1:B:337:THR:HA	1.62	0.81
1:B:59:GLY:HA3	1:B:64:ASN:HB3	1.63	0.81
1:D:289:LEU:HB3	1:D:290:PRO:HD2	1.60	0.81
1:D:294:ALA:H	1:D:337:THR:HG21	1.45	0.81
1:C:72:ILE:HG12	1:C:118:MSE:HE1	1.63	0.81
1:A:89:ASN:O	1:A:93:GLU:HG2	1.82	0.79
1:B:280:PHE:HE1	1:B:296:MSE:HE2	1.47	0.78
1:D:195:LEU:HD23	2:D:348:HOH:O	1.83	0.78
1:C:114:VAL:CG1	1:C:118:MSE:HE2	2.14	0.78
1:D:338:LEU:HD23	1:D:339:LYS:N	1.99	0.78
1:C:137:ALA:HA	1:C:140:MSE:CE	2.15	0.77
1:B:30:ARG:HH22	1:B:58:ALA:HB2	1.48	0.77
1:C:120:SER:HB2	1:D:56:LYS:HG2	1.67	0.77
1:D:322:ILE:HG22	1:D:323:GLU:N	1.99	0.77
2:A:376:HOH:O	1:D:221:GLU:HG3	1.86	0.75
1:D:40:ARG:NH1	1:D:44:LEU:HD11	2.02	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:VAL:HB	1:B:100:VAL:CG1	2.17	0.74
1:A:255:MSE:HA	1:A:255:MSE:CE	2.17	0.74
1:A:15:SER:O	1:A:19:GLY:N	2.21	0.74
1:D:32:ARG:HD3	1:D:32:ARG:N	2.04	0.73
1:B:78:ASN:O	1:B:79:MSE:HG2	1.90	0.72
1:A:55:ASN:ND2	1:B:61:LEU:H	1.87	0.71
1:C:208:GLU:O	1:C:212:LYS:HG2	1.90	0.71
1:A:35:VAL:HG13	1:A:39:THR:HB	1.73	0.71
1:B:89:ASN:O	1:B:93:GLU:HG2	1.91	0.71
1:C:62:ALA:HA	1:C:65:ARG:HH11	1.54	0.71
1:A:35:VAL:CG1	1:A:39:THR:HB	2.21	0.70
1:C:145:GLY:HA3	1:D:8:PRO:HB3	1.71	0.70
1:D:33:GLY:O	1:D:35:VAL:HG23	1.91	0.70
1:C:100:VAL:HB	1:D:100:VAL:HG22	1.72	0.70
1:A:200:ARG:O	1:A:204:GLU:HG3	1.90	0.70
1:C:79:MSE:HE1	2:D:364:HOH:O	1.91	0.70
1:A:27:ARG:NH1	1:A:33:GLY:HA3	2.06	0.70
1:D:61:LEU:HA	1:D:122:ARG:NE	2.07	0.69
1:A:255:MSE:HE1	1:A:285:LEU:HD13	1.73	0.69
1:D:286:LEU:HD11	1:D:293:LEU:HD23	1.74	0.69
1:D:293:LEU:HD12	1:D:337:THR:HG21	1.74	0.69
1:C:95:THR:HG22	1:C:97:LEU:N	2.05	0.68
1:D:10:THR:O	1:D:12:ARG:N	2.27	0.68
1:A:186:ILE:HG22	1:A:187:GLY:N	2.08	0.68
1:B:185:ARG:HB3	1:B:218:GLU:OE1	1.94	0.68
1:D:24:THR:O	1:D:28:VAL:HG23	1.93	0.68
1:A:11:LEU:C	1:A:13:ASP:H	2.02	0.68
1:A:27:ARG:HD3	1:A:32:ARG:HD3	1.75	0.67
1:D:29:LEU:O	1:D:30:ARG:HG2	1.94	0.67
1:C:57:ILE:HD13	1:D:121:TRP:O	1.95	0.67
1:D:21:SER:O	1:D:25:VAL:HG23	1.94	0.67
1:C:277:LEU:HB3	2:C:346:HOH:O	1.93	0.67
1:D:115:LEU:HD23	1:D:140:MSE:CE	2.24	0.67
2:A:367:HOH:O	1:B:96:GLU:HG3	1.95	0.67
1:B:170:ALA:HB2	1:B:296:MSE:HE1	1.74	0.67
1:D:338:LEU:HD23	1:D:339:LYS:H	1.57	0.67
1:D:221:GLU:HB3	1:D:235:MSE:HE2	1.77	0.66
1:D:57:ILE:HG22	1:D:66:VAL:HG11	1.78	0.66
1:C:95:THR:HG22	1:C:96:GLU:N	2.11	0.66
1:C:57:ILE:HG23	1:D:121:TRP:HA	1.78	0.66
1:D:338:LEU:O	1:D:339:LYS:HB2	1.93	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:SER:OG	1:A:43:VAL:HG12	1.96	0.66
1:C:72:ILE:HG12	1:C:118:MSE:CE	2.26	0.66
1:D:322:ILE:CG2	1:D:323:GLU:H	2.09	0.65
1:C:219:ASP:OD2	1:C:239:MSE:HG3	1.97	0.65
1:A:185:ARG:HB3	1:A:218:GLU:OE1	1.96	0.64
1:D:59:GLY:HA2	1:D:64:ASN:HB3	1.78	0.64
1:C:238:ALA:O	1:C:241:GLU:HB3	1.98	0.64
1:B:35:VAL:HG12	1:B:39:THR:HG23	1.80	0.64
1:D:270:ASP:OD2	1:D:273:GLY:HA3	1.98	0.64
1:B:277:LEU:CD1	1:B:337:THR:HB	2.27	0.63
1:B:264:LEU:HG	1:B:269:ILE:HD11	1.78	0.63
1:A:187:GLY:HA2	1:A:219:ASP:O	1.98	0.63
1:C:191:THR:HG22	1:C:223:TYR:HB3	1.81	0.63
1:A:55:ASN:HD22	1:B:61:LEU:N	1.95	0.63
1:B:280:PHE:O	1:B:281:ASN:HB2	1.98	0.62
1:B:190:GLY:HA2	1:B:252:SER:HB2	1.79	0.62
1:C:76:LEU:HD13	1:D:76:LEU:HD22	1.81	0.62
1:B:46:ALA:O	1:B:49:GLU:HG2	1.98	0.62
1:A:92:LEU:HD12	1:A:99:PRO:HG3	1.81	0.62
1:D:271:ILE:HD12	1:D:277:LEU:HD11	1.82	0.62
1:C:280:PHE:O	1:C:281:ASN:HB2	1.98	0.62
1:C:157:LYS:HB2	1:C:157:LYS:NZ	2.15	0.61
1:B:36:SER:OG	1:B:39:THR:HG22	2.00	0.61
1:B:320:ALA:O	1:B:321:GLU:HB3	2.00	0.61
1:B:339:LYS:O	1:B:340:ARG:HB3	1.99	0.61
1:C:115:LEU:CD2	1:C:140:MSE:HE3	2.26	0.61
1:D:115:LEU:CD2	1:D:140:MSE:HE3	2.31	0.61
1:C:192:LYS:HG3	1:C:196:ASP:HB2	1.83	0.61
1:A:57:ILE:HD11	1:B:57:ILE:HG21	1.83	0.60
1:A:21:SER:O	1:A:22:GLU:HB2	2.00	0.60
1:C:121:TRP:HA	1:D:57:ILE:HG23	1.83	0.60
1:C:89:ASN:O	1:C:93:GLU:HG2	2.02	0.60
1:A:39:THR:HA	1:A:42:ARG:NE	2.16	0.60
1:A:86:THR:O	1:A:90:GLN:HG3	2.02	0.60
1:A:39:THR:O	1:A:42:ARG:HB3	2.02	0.60
1:B:132:GLU:HG2	1:B:197:TYR:CD2	2.37	0.60
1:B:112:GLU:HG3	1:B:140:MSE:CE	2.31	0.60
1:A:12:ARG:O	1:A:12:ARG:HG3	2.01	0.60
1:C:105:ASP:O	1:C:106:TYR:HB2	2.02	0.60
1:B:155:ASP:OD1	1:B:201:LYS:HD3	2.02	0.59
1:B:12:ARG:O	1:B:16:GLU:HG3	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:PHE:CE1	1:B:296:MSE:HE2	2.35	0.59
1:C:283:VAL:HA	1:C:300:ARG:NH1	2.11	0.59
1:A:322:ILE:HG13	1:A:323:GLU:H	1.67	0.59
1:D:30:ARG:O	1:D:31:ASN:HB2	2.02	0.59
1:D:119:LEU:HD12	1:D:140:MSE:HG2	1.84	0.59
1:C:60:ALA:H	1:C:64:ASN:HB2	1.67	0.59
1:C:221:GLU:O	1:C:221:GLU:HG3	2.01	0.59
1:A:27:ARG:HD2	1:A:33:GLY:H	1.66	0.59
1:C:95:THR:CG2	1:C:96:GLU:N	2.66	0.59
1:A:83:GLU:HB3	1:A:300:ARG:HD2	1.85	0.59
1:B:79:MSE:HG3	1:B:255:MSE:CE	2.33	0.59
1:C:27:ARG:HE	1:C:32:ARG:HH12	1.51	0.59
1:C:301:LEU:O	1:C:305:ARG:HG3	2.02	0.59
1:A:30:ARG:HH11	1:A:30:ARG:HG2	1.68	0.58
1:B:184:ARG:O	1:B:247:ASP:OD1	2.22	0.58
1:B:336:ASP:O	1:B:338:LEU:N	2.37	0.58
1:D:20:VAL:HG12	1:D:42:ARG:HH21	1.68	0.58
1:A:38:ALA:O	1:A:42:ARG:HB2	2.03	0.58
1:A:40:ARG:O	1:A:44:LEU:HG	2.03	0.58
1:A:92:LEU:CD1	1:A:99:PRO:HG3	2.33	0.58
1:D:27:ARG:HG3	1:D:32:ARG:HH21	1.68	0.58
1:A:40:ARG:C	1:A:42:ARG:H	2.10	0.58
1:A:280:PHE:O	1:A:281:ASN:HB2	2.03	0.58
1:C:93:GLU:OE2	1:D:110:LYS:HE3	2.04	0.58
1:D:86:THR:O	1:D:90:GLN:HG3	2.04	0.58
1:B:59:GLY:CA	1:B:64:ASN:HB3	2.32	0.58
1:B:174:MSE:HG2	1:B:250:TYR:CE1	2.40	0.57
1:D:28:VAL:HG11	1:D:43:VAL:HB	1.86	0.57
1:D:294:ALA:N	1:D:337:THR:CG2	2.64	0.57
1:D:305:ARG:O	1:D:309:GLU:HG3	2.04	0.57
1:B:105:ASP:O	1:B:106:TYR:HB2	2.03	0.57
1:D:39:THR:O	1:D:43:VAL:HG23	2.05	0.57
1:A:9:LEU:C	1:A:9:LEU:HD23	2.29	0.57
1:A:68:LEU:HD22	1:A:123:PRO:HB3	1.86	0.57
1:A:83:GLU:OE1	1:A:83:GLU:HA	2.03	0.57
1:A:24:THR:O	1:A:26:SER:N	2.38	0.56
1:B:124:SER:HB3	1:B:315:LEU:HD21	1.86	0.56
1:C:80:VAL:HA	1:C:300:ARG:HD2	1.87	0.56
1:A:186:ILE:O	1:A:248:PHE:O	2.22	0.56
1:B:86:THR:O	1:B:90:GLN:HG3	2.05	0.56
1:C:100:VAL:HB	1:D:100:VAL:CG2	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:ARG:NH2	1:C:266:GLU:OE1	2.38	0.56
1:A:56:LYS:HB2	1:B:120:SER:HB2	1.88	0.56
1:C:187:GLY:HA2	1:C:219:ASP:O	2.06	0.56
1:D:280:PHE:O	1:D:281:ASN:HB2	2.05	0.56
1:D:20:VAL:HG23	1:D:25:VAL:HG22	1.89	0.56
1:A:122:ARG:HH21	1:B:55:ASN:HB2	1.71	0.55
1:B:320:ALA:O	1:B:321:GLU:CB	2.53	0.55
1:D:57:ILE:HG13	1:D:57:ILE:O	2.05	0.55
1:D:134:SER:O	1:D:135:GLU:HB3	2.06	0.55
1:A:186:ILE:HD12	1:A:186:ILE:N	2.22	0.55
1:C:57:ILE:CG2	1:D:121:TRP:HA	2.36	0.55
1:C:114:VAL:HG12	1:C:118:MSE:CE	2.34	0.55
1:D:30:ARG:HE	1:D:30:ARG:HA	1.72	0.55
1:A:27:ARG:CD	1:A:32:ARG:HB3	2.37	0.55
1:C:193:MSE:HE2	1:C:222:PHE:CE2	2.41	0.55
1:A:162:MSE:HE2	1:A:325:ARG:HB3	1.88	0.54
1:A:221:GLU:HG3	1:A:235:MSE:HE3	1.87	0.54
1:C:193:MSE:HE3	1:C:200:ARG:HG2	1.89	0.54
1:B:264:LEU:CG	1:B:269:ILE:HD11	2.37	0.54
1:D:313:LYS:HE3	1:D:322:ILE:HG13	1.89	0.54
1:A:76:LEU:HD13	1:B:76:LEU:HD22	1.88	0.54
1:C:61:LEU:HG	1:C:122:ARG:HH21	1.72	0.54
1:A:11:LEU:C	1:A:13:ASP:N	2.66	0.54
1:B:172:ARG:HG2	1:B:205:GLY:O	2.08	0.54
1:D:79:MSE:HE2	1:D:255:MSE:HE3	1.89	0.54
1:A:27:ARG:HD3	1:A:32:ARG:HB3	1.90	0.54
1:B:57:ILE:HG13	1:B:57:ILE:O	2.08	0.54
1:A:322:ILE:HG13	1:A:323:GLU:N	2.23	0.53
1:C:285:LEU:HD11	1:D:285:LEU:HD11	1.90	0.53
1:B:79:MSE:HG3	1:B:255:MSE:HE1	1.91	0.53
1:C:223:TYR:CE2	1:C:231:LYS:HE2	2.43	0.53
1:A:68:LEU:C	1:A:68:LEU:HD23	2.34	0.53
1:B:89:ASN:ND2	1:B:99:PRO:HG3	2.24	0.53
1:A:18:SER:O	1:A:20:VAL:HG23	2.09	0.53
1:B:264:LEU:HG	1:B:269:ILE:CD1	2.39	0.53
1:D:148:VAL:O	1:D:161:ALA:HB3	2.08	0.53
1:B:30:ARG:NH2	1:B:58:ALA:HB2	2.22	0.52
1:A:147:PRO:HB3	1:A:314:ARG:NH1	2.25	0.52
1:B:170:ALA:CB	1:B:296:MSE:HE1	2.39	0.52
1:B:325:ARG:HD2	2:B:366:HOH:O	2.08	0.52
1:C:14:VAL:O	1:C:18:SER:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:ASN:ND2	1:B:117:GLU:OE1	2.42	0.52
1:A:237:GLN:HG3	1:A:238:ALA:N	2.24	0.52
1:A:204:GLU:O	1:A:208:GLU:HG3	2.09	0.52
1:B:194:PRO:HA	1:B:200:ARG:HD2	1.91	0.52
1:C:184:ARG:O	1:C:247:ASP:OD1	2.28	0.52
1:D:28:VAL:HG13	1:D:40:ARG:HG3	1.92	0.52
1:A:27:ARG:CZ	1:A:33:GLY:HA3	2.38	0.52
1:B:174:MSE:O	1:B:178:ILE:HG13	2.10	0.52
1:C:169:ARG:NH1	1:C:169:ARG:HG3	2.24	0.52
1:A:52:TYR:OH	1:B:61:LEU:HD21	2.09	0.52
1:C:81:PHE:O	1:C:85:LEU:HG	2.10	0.52
1:D:18:SER:OG	1:D:20:VAL:HG22	2.10	0.52
1:D:30:ARG:HB3	1:D:32:ARG:HD2	1.91	0.52
1:A:255:MSE:HE1	1:A:285:LEU:HD22	1.92	0.51
1:C:86:THR:O	1:C:90:GLN:HG3	2.10	0.51
1:D:178:ILE:CG2	1:D:183:TYR:HB2	2.40	0.51
1:B:112:GLU:HG3	1:B:140:MSE:HE1	1.91	0.51
1:C:72:ILE:HD11	1:C:118:MSE:HE3	1.92	0.51
1:D:79:MSE:HE2	1:D:255:MSE:SE	2.61	0.51
1:B:178:ILE:HD12	1:B:206:PHE:HE1	1.76	0.51
1:C:220:ARG:HG2	1:C:220:ARG:HH11	1.76	0.51
1:A:79:MSE:HE1	1:A:285:LEU:HD13	1.91	0.51
1:C:72:ILE:CG1	1:C:118:MSE:HE1	2.39	0.51
1:C:183:TYR:OH	1:C:272:PRO:HB2	2.11	0.51
1:A:276:GLY:C	1:A:277:LEU:HD12	2.36	0.50
1:D:77:SER:C	1:D:78:ASN:O	2.52	0.50
1:B:190:GLY:HA2	1:B:252:SER:CB	2.40	0.50
1:B:249:LEU:CG	1:B:275:ILE:HD11	2.29	0.50
1:D:160:ASP:O	1:D:161:ALA:O	2.30	0.50
1:B:219:ASP:OD2	1:B:242:ARG:HD2	2.12	0.50
1:C:8:PRO:HB2	1:D:62:ALA:HB2	1.94	0.50
1:C:306:LYS:HE3	1:C:326:ILE:CD1	2.42	0.50
1:D:115:LEU:CG	1:D:140:MSE:HE3	2.42	0.50
1:C:255:MSE:SE	1:C:285:LEU:HD13	2.61	0.50
1:A:36:SER:O	1:A:40:ARG:HG2	2.12	0.50
1:A:198:ARG:HG3	1:A:198:ARG:HH11	1.77	0.50
1:C:27:ARG:HE	1:C:32:ARG:NH1	2.10	0.50
1:B:49:GLU:HG3	1:B:50:LEU:N	2.27	0.49
1:C:83:GLU:HG2	1:C:300:ARG:HE	1.77	0.49
1:C:109:GLU:CD	1:C:109:GLU:H	2.20	0.49
1:B:317:ASP:HB3	1:B:320:ALA:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:VAL:O	1:D:160:ASP:O	2.29	0.49
1:D:193:MSE:HE3	1:D:200:ARG:HA	1.95	0.49
1:A:280:PHE:O	1:A:281:ASN:CB	2.60	0.49
1:C:220:ARG:HD2	1:C:220:ARG:O	2.12	0.49
1:D:6:LYS:NZ	1:D:6:LYS:HB3	2.26	0.49
1:A:319:GLU:O	1:A:320:ALA:C	2.56	0.49
1:C:192:LYS:O	1:C:192:LYS:HD3	2.13	0.49
1:C:193:MSE:HA	1:C:195:LEU:H	1.77	0.49
1:D:292:LYS:O	1:D:335:GLY:HA3	2.13	0.49
1:C:95:THR:CG2	1:C:96:GLU:H	2.26	0.49
1:B:277:LEU:HD13	1:B:293:LEU:CD1	2.42	0.49
1:C:277:LEU:HD23	1:C:277:LEU:N	2.28	0.49
1:D:285:LEU:C	1:D:285:LEU:HD23	2.38	0.49
1:C:85:LEU:HD13	1:C:101:VAL:HG21	1.95	0.48
1:D:105:ASP:O	1:D:106:TYR:HB2	2.13	0.48
1:D:280:PHE:O	1:D:281:ASN:CB	2.60	0.48
1:A:231:LYS:NZ	1:A:234:GLU:OE1	2.44	0.48
1:D:257:ALA:HB1	1:D:293:LEU:HD21	1.96	0.48
1:C:28:VAL:C	1:C:30:ARG:H	2.21	0.48
1:A:85:LEU:HD22	1:A:101:VAL:HG21	1.95	0.48
1:A:128:ILE:O	1:A:150:GLU:HA	2.13	0.48
1:B:338:LEU:C	1:B:338:LEU:HD23	2.38	0.48
1:C:169:ARG:HG3	1:C:169:ARG:HH11	1.78	0.48
1:B:155:ASP:O	1:B:156:GLY:O	2.32	0.48
1:C:221:GLU:HB3	1:C:242:ARG:HH22	1.77	0.48
1:C:229:LEU:HD12	1:C:255:MSE:HG3	1.96	0.48
1:C:251:TYR:HB3	1:C:256:ILE:HG13	1.96	0.48
1:C:261:LEU:O	1:C:265:LEU:HG	2.14	0.48
1:D:296:MSE:HE3	1:D:332:ILE:HG13	1.94	0.48
1:A:93:GLU:O	1:A:95:THR:N	2.46	0.47
1:D:20:VAL:HG23	1:D:25:VAL:CG2	2.44	0.47
1:D:221:GLU:HB3	1:D:235:MSE:CE	2.44	0.47
1:C:191:THR:HA	1:C:223:TYR:O	2.14	0.47
1:D:184:ARG:O	1:D:247:ASP:CG	2.57	0.47
1:A:160:ASP:CG	1:A:314:ARG:HH22	2.21	0.47
1:A:24:THR:O	1:A:25:VAL:C	2.57	0.47
1:B:218:GLU:HB3	1:B:246:LEU:HD21	1.96	0.47
1:C:11:LEU:HD22	1:C:25:VAL:HG12	1.96	0.47
1:A:255:MSE:HE1	1:A:285:LEU:CD1	2.44	0.47
1:B:338:LEU:O	1:B:339:LYS:C	2.58	0.47
1:D:322:ILE:CG2	1:D:323:GLU:N	2.68	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:ASN:C	1:A:93:GLU:HG2	2.40	0.47
1:A:121:TRP:CD1	1:B:66:VAL:HG21	2.50	0.47
1:C:83:GLU:HB3	1:C:300:ARG:HG2	1.96	0.47
1:A:174:MSE:HG2	1:A:250:TYR:CE1	2.49	0.47
1:B:62:ALA:O	1:B:63:SER:CB	2.62	0.47
1:C:223:TYR:CD2	1:C:231:LYS:HE2	2.50	0.47
1:D:229:LEU:HD12	1:D:255:MSE:HG3	1.97	0.47
1:A:105:ASP:O	1:A:106:TYR:HB2	2.14	0.46
1:C:78:ASN:O	1:C:79:MSE:HE2	2.15	0.46
1:C:221:GLU:OE2	1:C:235:MSE:HG2	2.15	0.46
1:C:193:MSE:HA	1:C:195:LEU:N	2.31	0.46
1:C:307:ALA:O	1:C:311:ILE:HG12	2.14	0.46
1:A:311:ILE:O	1:A:315:LEU:HD13	2.15	0.46
1:C:272:PRO:HA	1:C:275:ILE:O	2.15	0.46
1:D:82:PRO:HG2	2:D:364:HOH:O	2.16	0.46
1:B:254:ASP:OD2	1:B:282:ASN:N	2.48	0.46
1:D:118:MSE:O	1:D:123:PRO:HD3	2.15	0.46
1:A:27:ARG:HD2	1:A:33:GLY:N	2.31	0.46
1:B:149:VAL:HG22	1:B:161:ALA:HB3	1.98	0.46
1:C:314:ARG:NH2	1:C:324:THR:HG21	2.31	0.46
1:B:61:LEU:HG	1:B:122:ARG:HH21	1.80	0.46
1:C:32:ARG:HG2	1:C:33:GLY:N	2.30	0.46
1:C:113:LYS:O	1:C:117:GLU:HG2	2.16	0.46
1:D:115:LEU:HB3	1:D:140:MSE:HE3	1.98	0.46
1:D:179:LEU:HD21	1:D:215:VAL:HG11	1.97	0.46
1:A:198:ARG:HG3	1:A:198:ARG:NH1	2.31	0.46
1:C:271:ILE:CG2	1:C:337:THR:HG22	2.46	0.46
1:A:192:LYS:HB2	1:A:192:LYS:NZ	2.31	0.46
1:D:40:ARG:HH12	1:D:44:LEU:HD11	1.77	0.46
1:D:317:ASP:OD2	1:D:320:ALA:HB2	2.15	0.46
1:A:149:VAL:HG22	1:A:161:ALA:HB3	1.98	0.46
1:B:61:LEU:HB3	1:B:62:ALA:H	1.50	0.46
1:B:184:ARG:C	1:B:185:ARG:HG3	2.40	0.46
1:D:20:VAL:HG21	1:D:43:VAL:CG2	2.37	0.46
1:D:272:PRO:HB3	1:D:337:THR:O	2.16	0.46
1:C:231:LYS:HG3	1:C:235:MSE:HE3	1.97	0.46
1:A:83:GLU:HB2	1:A:300:ARG:NH1	2.31	0.45
1:A:338:LEU:C	1:A:338:LEU:HD23	2.41	0.45
1:C:233:ARG:HH22	1:C:266:GLU:CD	2.24	0.45
1:A:40:ARG:C	1:A:42:ARG:N	2.74	0.45
1:C:55:ASN:ND2	1:D:61:LEU:H	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:157:LYS:HB3	1:C:325:ARG:NH1	2.31	0.45
1:D:298:ALA:HB3	1:D:300:ARG:HG3	1.98	0.45
1:A:189:MSE:HG2	1:A:221:GLU:HB3	1.98	0.45
1:C:121:TRP:CZ3	1:D:68:LEU:HD12	2.52	0.45
1:D:79:MSE:O	1:D:300:ARG:NH1	2.49	0.45
1:B:160:ASP:CG	1:B:314:ARG:HH12	2.25	0.45
1:C:280:PHE:O	1:C:281:ASN:CB	2.63	0.45
1:C:338:LEU:HD23	1:C:338:LEU:C	2.42	0.45
1:B:85:LEU:HD22	1:B:101:VAL:HG21	1.99	0.45
1:B:280:PHE:O	1:B:281:ASN:CB	2.64	0.45
1:C:72:ILE:CG1	1:C:118:MSE:CE	2.95	0.45
1:C:80:VAL:O	1:C:300:ARG:HG2	2.17	0.45
1:C:188:PHE:HB3	1:C:220:ARG:HB3	2.00	0.44
1:A:165:ILE:HD12	1:A:166:SER:C	2.42	0.44
1:A:184:ARG:O	1:A:247:ASP:OD1	2.35	0.44
1:A:317:ASP:O	1:A:320:ALA:N	2.47	0.44
1:C:157:LYS:HB2	1:C:157:LYS:HZ3	1.82	0.44
1:C:239:MSE:C	1:C:241:GLU:H	2.26	0.44
1:D:293:LEU:HA	1:D:337:THR:HG21	1.99	0.44
1:A:29:LEU:HD12	1:A:44:LEU:CD2	2.47	0.44
1:D:169:ARG:O	1:D:173:GLU:HB2	2.18	0.44
1:A:36:SER:C	1:A:38:ALA:H	2.26	0.44
1:C:30:ARG:O	1:C:31:ASN:HB2	2.16	0.44
1:D:29:LEU:HD22	1:D:54:PRO:HB3	2.00	0.44
1:D:264:LEU:HD23	1:D:271:ILE:HD11	1.98	0.44
1:B:129:ALA:HA	1:B:151:ILE:HG13	1.99	0.44
1:C:320:ALA:O	1:C:321:GLU:HB3	2.18	0.43
1:C:185:ARG:HB3	1:C:218:GLU:OE1	2.17	0.43
1:A:69:VAL:HG23	1:A:97:LEU:HB3	1.99	0.43
1:A:318:PRO:C	1:A:320:ALA:H	2.25	0.43
1:C:156:GLY:O	1:C:157:LYS:C	2.60	0.43
1:A:31:ASN:HA	1:A:40:ARG:NH2	2.24	0.43
1:D:231:LYS:HD2	1:D:231:LYS:HA	1.86	0.43
1:D:272:PRO:HD3	1:D:336:ASP:O	2.18	0.43
1:B:61:LEU:HA	1:B:122:ARG:HE	1.82	0.43
1:B:315:LEU:HD12	1:B:315:LEU:HA	1.82	0.43
1:D:97:LEU:HD11	1:D:315:LEU:HD23	2.00	0.43
1:C:229:LEU:HD11	1:C:255:MSE:SE	2.69	0.43
1:D:134:SER:O	1:D:135:GLU:CB	2.67	0.43
1:D:184:ARG:O	1:D:247:ASP:OD1	2.36	0.43
1:A:20:VAL:CG1	1:A:24:THR:OG1	2.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:59:GLY:HA3	1:C:64:ASN:HB3	2.00	0.43
1:A:27:ARG:HD3	1:A:32:ARG:CD	2.44	0.43
1:A:83:GLU:HB2	1:A:300:ARG:HH11	1.84	0.43
1:A:302:GLU:H	1:A:302:GLU:CD	2.26	0.43
1:B:292:LYS:O	1:B:335:GLY:O	2.37	0.43
1:C:183:TYR:CD2	1:C:248:PHE:HB2	2.54	0.43
1:D:192:LYS:HA	1:D:192:LYS:HD3	1.76	0.43
1:B:35:VAL:CG1	1:B:39:THR:HG23	2.47	0.43
1:C:36:SER:C	1:C:38:ALA:H	2.27	0.43
1:C:61:LEU:HB3	1:C:62:ALA:H	1.53	0.43
1:C:171:GLY:O	1:C:206:PHE:HA	2.18	0.43
1:C:271:ILE:HG21	1:C:337:THR:HG22	2.01	0.43
1:A:160:ASP:OD2	1:A:314:ARG:NH2	2.52	0.43
1:B:257:ALA:HB1	1:B:293:LEU:HD21	2.01	0.43
1:B:295:THR:OG1	1:B:296:MSE:N	2.51	0.43
1:B:59:GLY:HA3	1:B:64:ASN:HD22	1.84	0.42
1:C:272:PRO:O	1:C:273:GLY:C	2.62	0.42
1:D:178:ILE:HG22	1:D:183:TYR:HB2	2.01	0.42
1:D:257:ALA:CB	1:D:293:LEU:HD21	2.49	0.42
1:A:118:MSE:O	1:A:123:PRO:HD3	2.19	0.42
1:C:191:THR:C	1:C:193:MSE:H	2.27	0.42
1:D:61:LEU:HB3	1:D:62:ALA:H	1.67	0.42
1:A:27:ARG:HD3	1:A:32:ARG:CG	2.49	0.42
1:C:254:ASP:OD2	1:C:282:ASN:N	2.52	0.42
1:D:294:ALA:N	1:D:337:THR:HG22	2.35	0.42
1:A:122:ARG:NH2	1:B:55:ASN:HB2	2.34	0.42
1:A:255:MSE:CE	1:A:285:LEU:HD22	2.49	0.42
1:B:107:LEU:HA	1:B:108:PRO:HD3	1.91	0.42
1:D:144:ALA:C	1:D:146:ILE:H	2.27	0.42
1:A:39:THR:HA	1:A:42:ARG:CZ	2.49	0.42
1:A:326:ILE:HD12	1:A:326:ILE:N	2.34	0.42
1:C:104:THR:O	1:C:110:LYS:HD2	2.20	0.42
1:C:178:ILE:HD12	1:C:206:PHE:HE1	1.85	0.42
1:A:237:GLN:HB2	1:A:263:TYR:CZ	2.54	0.42
1:C:115:LEU:HD23	1:C:140:MSE:CE	2.29	0.42
1:C:128:ILE:O	1:C:150:GLU:HA	2.20	0.42
1:D:109:GLU:HG2	1:D:110:LYS:N	2.35	0.42
1:D:263:TYR:O	1:D:267:GLN:HG2	2.19	0.42
1:C:145:GLY:CA	1:D:8:PRO:HB3	2.46	0.42
1:D:206:PHE:CE2	1:D:217:ILE:HD11	2.55	0.42
1:D:293:LEU:HA	1:D:337:THR:CG2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:PRO:O	1:A:9:LEU:O	2.37	0.42
1:A:65:ARG:HD3	1:A:123:PRO:O	2.20	0.42
1:C:12:ARG:O	1:C:16:GLU:HG3	2.20	0.42
1:D:185:ARG:HD3	1:D:245:ASP:O	2.20	0.42
1:A:11:LEU:O	1:A:13:ASP:N	2.52	0.42
1:A:184:ARG:O	1:A:247:ASP:CG	2.63	0.42
1:B:29:LEU:C	1:B:31:ASN:H	2.28	0.42
1:B:118:MSE:O	1:B:123:PRO:HD3	2.20	0.42
1:C:254:ASP:HB3	1:C:286:LEU:HD11	2.02	0.42
1:D:20:VAL:CG1	1:D:42:ARG:HH21	2.31	0.42
1:D:276:GLY:C	1:D:277:LEU:HD12	2.45	0.42
1:C:73:ILE:HB	1:C:74:PRO:HD2	2.01	0.42
1:C:131:LEU:HD22	1:C:158:PRO:CD	2.50	0.42
1:D:30:ARG:HA	1:D:30:ARG:NE	2.33	0.42
1:A:158:PRO:HB3	1:A:162:MSE:HB2	2.02	0.41
1:D:147:PRO:HG2	1:D:315:LEU:CD1	2.50	0.41
1:A:47:ALA:HB1	1:A:52:TYR:HB3	2.03	0.41
1:A:124:SER:HB3	1:A:315:LEU:HD21	2.02	0.41
1:A:57:ILE:HD11	1:B:57:ILE:CG2	2.48	0.41
1:B:89:ASN:ND2	1:B:99:PRO:CG	2.83	0.41
1:A:28:VAL:HG11	1:A:40:ARG:HA	2.02	0.41
1:B:94:ASP:OD2	1:C:331:LYS:CB	2.68	0.41
1:B:184:ARG:O	1:B:185:ARG:CB	2.67	0.41
1:A:262:LEU:O	1:A:266:GLU:HB2	2.20	0.41
1:B:128:ILE:O	1:B:150:GLU:HA	2.19	0.41
1:C:239:MSE:C	1:C:241:GLU:N	2.78	0.41
1:C:285:LEU:HG	1:D:229:LEU:HD22	2.02	0.41
1:D:79:MSE:CE	1:D:255:MSE:HE3	2.49	0.41
1:D:271:ILE:HD12	1:D:277:LEU:CD1	2.49	0.41
1:C:306:LYS:HD2	1:C:306:LYS:HA	1.84	0.41
1:A:55:ASN:ND2	1:B:61:LEU:N	2.60	0.41
1:A:160:ASP:OD2	1:A:160:ASP:C	2.64	0.41
1:C:231:LYS:HD2	1:C:231:LYS:HA	1.91	0.41
1:D:289:LEU:HB3	1:D:290:PRO:CD	2.42	0.41
1:D:296:MSE:HE3	1:D:332:ILE:CG1	2.50	0.41
1:C:100:VAL:CG2	1:D:100:VAL:HG21	2.51	0.41
1:A:27:ARG:HD2	1:A:32:ARG:HB3	2.02	0.41
1:B:66:VAL:HG22	1:B:67:ASN:N	2.36	0.41
1:C:68:LEU:HD23	1:C:69:VAL:N	2.35	0.41
1:D:79:MSE:HE2	1:D:255:MSE:CE	2.51	0.41
1:B:223:TYR:CE1	1:B:231:LYS:HE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:57:ILE:HD11	1:D:122:ARG:HB2	2.03	0.40
1:C:193:MSE:O	1:C:224:SER:HA	2.21	0.40
1:D:106:TYR:HB3	1:D:195:LEU:HB3	2.01	0.40
1:D:128:ILE:O	1:D:150:GLU:HA	2.22	0.40
1:D:223:TYR:CD2	1:D:231:LYS:HE2	2.56	0.40
1:A:167:HIS:O	1:A:202:ARG:HA	2.21	0.40
1:A:192:LYS:HD3	1:A:192:LYS:HA	1.76	0.40
1:A:264:LEU:HD22	1:A:277:LEU:HD21	2.03	0.40
1:B:28:VAL:HG22	1:B:40:ARG:HG2	2.03	0.40
1:C:57:ILE:HG22	1:C:66:VAL:HG21	2.03	0.40
1:C:119:LEU:HD23	1:C:119:LEU:HA	1.94	0.40
1:C:302:GLU:O	1:C:306:LYS:HB2	2.22	0.40
1:C:97:LEU:HD11	1:C:315:LEU:HD12	2.04	0.40
1:C:283:VAL:O	1:C:286:LEU:HB2	2.21	0.40
1:C:292:LYS:HE2	1:C:336:ASP:OD2	2.22	0.40
1:D:223:TYR:CG	1:D:231:LYS:HE2	2.55	0.40
1:D:320:ALA:O	1:D:321:GLU:O	2.40	0.40
1:A:83:GLU:CB	1:A:300:ARG:HD2	2.50	0.40
1:A:192:LYS:HB2	1:A:192:LYS:HZ2	1.86	0.40
1:C:107:LEU:HD12	1:C:110:LYS:HE3	2.04	0.40
1:C:284:GLU:CD	1:C:284:GLU:C	2.89	0.40
1:D:27:ARG:HH21	1:D:33:GLY:HA3	1.87	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	331/344 (96%)	285 (86%)	32 (10%)	14 (4%)	2 1
1	B	331/344 (96%)	303 (92%)	17 (5%)	11 (3%)	3 1
1	C	331/344 (96%)	293 (88%)	26 (8%)	12 (4%)	2 1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	D	333/344 (97%)	298 (90%)	24 (7%)	11 (3%)	3 1
All	All	1326/1376 (96%)	1179 (89%)	99 (8%)	48 (4%)	2 1

All (48) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	LEU
1	A	22	GLU
1	A	25	VAL
1	A	61	LEU
1	B	61	LEU
1	B	79	MSE
1	B	155	ASP
1	B	156	GLY
1	B	281	ASN
1	B	337	THR
1	C	61	LEU
1	C	158	PRO
1	C	281	ASN
1	D	11	LEU
1	D	161	ALA
1	D	321	GLU
1	A	24	THR
1	A	186	ILE
1	A	281	ASN
1	A	320	ALA
1	B	63	SER
1	B	339	LYS
1	C	29	LEU
1	C	59	GLY
1	C	222	PHE
1	D	30	ARG
1	D	32	ARG
1	D	63	SER
1	D	281	ASN
1	A	12	ARG
1	A	35	VAL
1	A	58	ALA
1	A	95	THR
1	B	78	ASN
1	B	321	GLU
1	C	35	VAL

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Mol	Chain	Res	Type
1	C	152	MSE
1	C	159	VAL
1	C	321	GLU
1	C	338	LEU
1	D	78	ASN
1	A	60	ALA
1	D	35	VAL
1	D	36	SER
1	A	21	SER
1	B	152	MSE
1	D	57	ILE
1	C	156	GLY

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	269/264 (102%)	257 (96%)	12 (4%)	24	32
1	B	269/264 (102%)	259 (96%)	10 (4%)	30	39
1	C	269/264 (102%)	259 (96%)	10 (4%)	30	39
1	D	271/264 (103%)	259 (96%)	12 (4%)	25	33
All	All	1078/1056 (102%)	1034 (96%)	44 (4%)	27	35

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ARG
1	A	23	MSE
1	A	24	THR
1	A	27	ARG
1	A	29	LEU
1	A	56	LYS
1	A	192	LYS
1	A	237	GLN
1	A	255	MSE

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Mol	Chain	Res	Type
1	A	286	LEU
1	A	306	LYS
1	A	321	GLU
1	B	11	LEU
1	B	29	LEU
1	B	39	THR
1	B	78	ASN
1	B	100	VAL
1	B	204	GLU
1	B	277	LEU
1	B	284	GLU
1	B	286	LEU
1	B	315	LEU
1	C	109	GLU
1	C	157	LYS
1	C	169	ARG
1	C	220	ARG
1	C	234	GLU
1	C	255	MSE
1	C	277	LEU
1	C	286	LEU
1	C	300	ARG
1	C	336	ASP
1	D	32	ARG
1	D	78	ASN
1	D	93	GLU
1	D	140	MSE
1	D	168	ARG
1	D	194	PRO
1	D	207	THR
1	D	220	ARG
1	D	277	LEU
1	D	285	LEU
1	D	315	LEU
1	D	329	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	ASN
1	A	287	GLN
1	B	31	ASN

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Mol	Chain	Res	Type
1	B	64	ASN
1	B	89	ASN
1	B	98	GLN
1	C	64	ASN
1	C	89	ASN
1	C	176	GLN
1	C	287	GLN
1	D	89	ASN
1	D	98	GLN
1	D	281	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](#)

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	320/344 (93%)	0.68	45 (14%) 6 7	22, 38, 107, 121	0
1	B	320/344 (93%)	0.56	24 (7%) 20 24	24, 43, 68, 88	0
1	C	320/344 (93%)	0.53	28 (8%) 15 19	19, 41, 75, 84	0
1	D	322/344 (93%)	0.41	34 (10%) 11 13	19, 34, 84, 113	0
All	All	1282/1376 (93%)	0.54	131 (10%) 12 14	19, 39, 83, 121	0

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	ALA	8.0
1	B	61	LEU	6.7
1	A	61	LEU	6.0
1	C	60	ALA	5.9
1	D	29	LEU	5.8
1	D	60	ALA	5.8
1	C	222	PHE	5.8
1	C	61	LEU	5.7
1	D	61	LEU	5.6
1	D	62	ALA	5.5
1	A	25	VAL	5.5
1	B	62	ALA	5.3
1	A	8	PRO	5.2
1	A	29	LEU	4.7
1	A	322	ILE	4.7
1	D	58	ALA	4.6
1	A	22	GLU	4.5
1	D	322	ILE	4.4
1	D	35	VAL	4.4
1	A	41	ALA	4.4
1	D	63	SER	4.3

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Mol	Chain	Res	Type	RSRZ
1	C	59	GLY	4.1
1	B	60	ALA	4.1
1	A	35	VAL	4.1
1	C	62	ALA	4.1
1	A	59	GLY	3.9
1	A	95	THR	3.9
1	C	63	SER	3.8
1	C	58	ALA	3.8
1	D	287	GLN	3.8
1	A	9	LEU	3.7
1	D	59	GLY	3.7
1	D	57	ILE	3.6
1	D	38	ALA	3.5
1	A	10	THR	3.3
1	D	39	THR	3.3
1	C	194	PRO	3.3
1	A	324	THR	3.3
1	B	63	SER	3.3
1	A	144	ALA	3.2
1	C	57	ILE	3.2
1	A	28	VAL	3.2
1	B	57	ILE	3.2
1	C	195	LEU	3.1
1	A	33	GLY	3.1
1	A	34	ASP	3.1
1	A	26	SER	3.1
1	C	223	TYR	3.0
1	B	58	ALA	3.0
1	D	28	VAL	3.0
1	A	21	SER	3.0
1	A	14	VAL	3.0
1	C	228	ALA	2.9
1	D	25	VAL	2.9
1	B	33	GLY	2.9
1	D	36	SER	2.9
1	A	20	VAL	2.9
1	A	63	SER	2.8
1	C	64	ASN	2.9
1	D	34	ASP	2.8
1	A	62	ALA	2.8
1	A	57	ILE	2.8
1	C	78	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	225	GLY	2.8
1	C	35	VAL	2.8
1	D	33	GLY	2.7
1	C	157	LYS	2.7
1	B	222	PHE	2.7
1	C	269	ILE	2.7
1	D	285	LEU	2.7
1	A	78	ASN	2.7
1	B	320	ALA	2.7
1	A	318	PRO	2.7
1	D	24	THR	2.6
1	C	158	PRO	2.6
1	B	339	LYS	2.6
1	D	284	GLU	2.6
1	A	30	ARG	2.6
1	A	39	THR	2.6
1	A	186	ILE	2.6
1	B	185	ARG	2.6
1	C	32	ARG	2.6
1	C	227	SER	2.6
1	C	321	GLU	2.5
1	A	17	ALA	2.5
1	A	24	THR	2.5
1	B	35	VAL	2.5
1	A	12	ARG	2.5
1	B	30	ARG	2.5
1	B	59	GLY	2.5
1	B	269	ILE	2.5
1	A	325	ARG	2.4
1	A	18	SER	2.4
1	D	31	ASN	2.4
1	A	42	ARG	2.4
1	D	26	SER	2.4
1	B	156	GLY	2.4
1	C	191	THR	2.4
1	A	326	ILE	2.4
1	A	11	LEU	2.3
1	B	273	GLY	2.3
1	A	38	ALA	2.3
1	A	315	LEU	2.3
1	D	40	ARG	2.3
1	B	31	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	241	GLU	2.2
1	D	11	LEU	2.2
1	B	64	ASN	2.2
1	D	324	THR	2.2
1	B	78	ASN	2.2
1	D	30	ARG	2.2
1	C	221	GLU	2.2
1	A	43	VAL	2.2
1	D	288	GLY	2.2
1	C	8	PRO	2.1
1	D	221	GLU	2.1
1	D	41	ALA	2.1
1	B	270	ASP	2.1
1	A	16	GLU	2.1
1	B	176	GLN	2.1
1	D	10	THR	2.1
1	A	13	ASP	2.1
1	B	49	GLU	2.0
1	D	144	ALA	2.0
1	D	320	ALA	2.0
1	A	44	LEU	2.0
1	D	37	ASP	2.0
1	B	340	ARG	2.0
1	A	19	GLY	2.0
1	C	39	THR	2.0
1	C	246	LEU	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.