



Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 02:10 am BST

PDB ID : 3U1V
Title : X-ray Structure of De Novo design cysteine esterase FR29, Northeast Structural Genomics Consortium Target OR52
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Deposited on : 2011-09-30
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

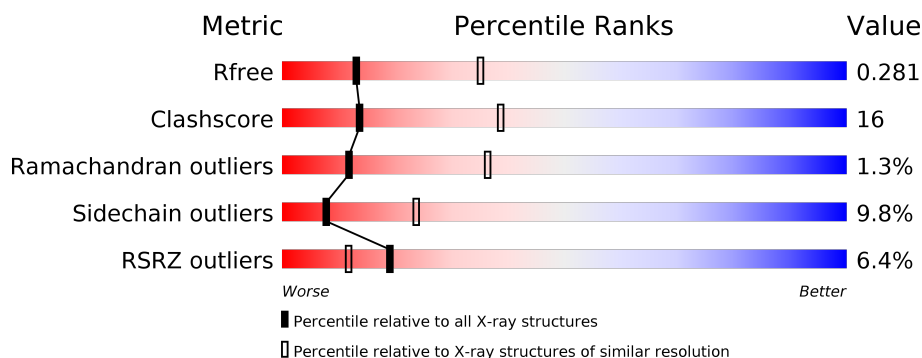
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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

Mol	Chain	Length	Quality of chain
1	A	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 67%, yellow 26%, orange 3%, grey 4%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 67% 26% • • </div> </div>
1	B	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 59%, yellow 34%, orange 3%, grey 4%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 59% 34% • • </div> </div>
1	C	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 17%, green 60%, yellow 34%, orange 3%, grey 4%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 17% 60% 34% • • </div> </div>
1	D	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 4%, green 48%, yellow 24%, orange 8%, grey 20%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 4% 48% 24% 8% 20% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9968 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 De Novo design cysteine esterase FR29.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	S	Se	0	0	0
			2597	1653	443	487	4	10			
1	B	327	Total	C	N	O	S	Se	0	0	0
			2597	1653	443	487	4	10			
1	C	327	Total	C	N	O	S	Se	0	0	0
			2597	1653	443	487	4	10			
1	D	269	Total	C	N	O	S	Se	0	0	0
			2145	1371	368	393	4	9			

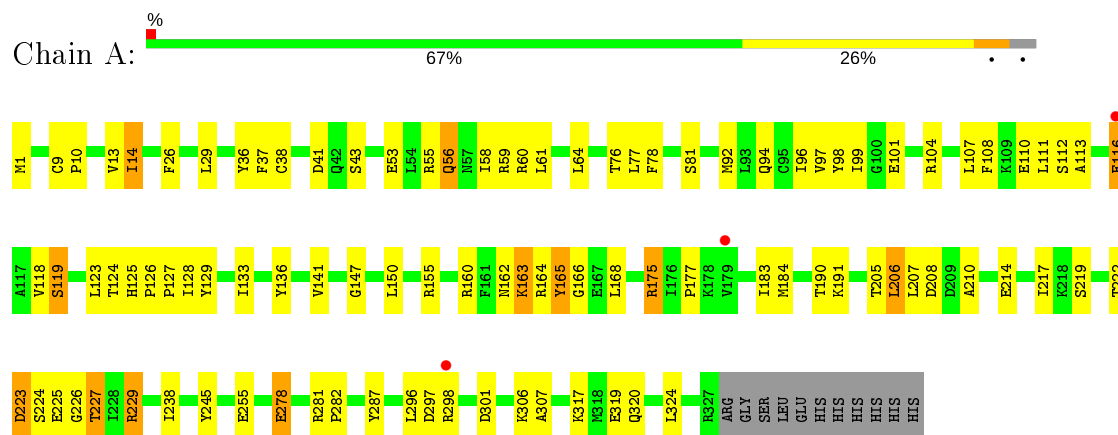
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	O	0	0
			9	9		
2	B	12	Total	O	0	0
			12	12		
2	C	6	Total	O	0	0
			6	6		
2	D	5	Total	O	0	0
			5	5		

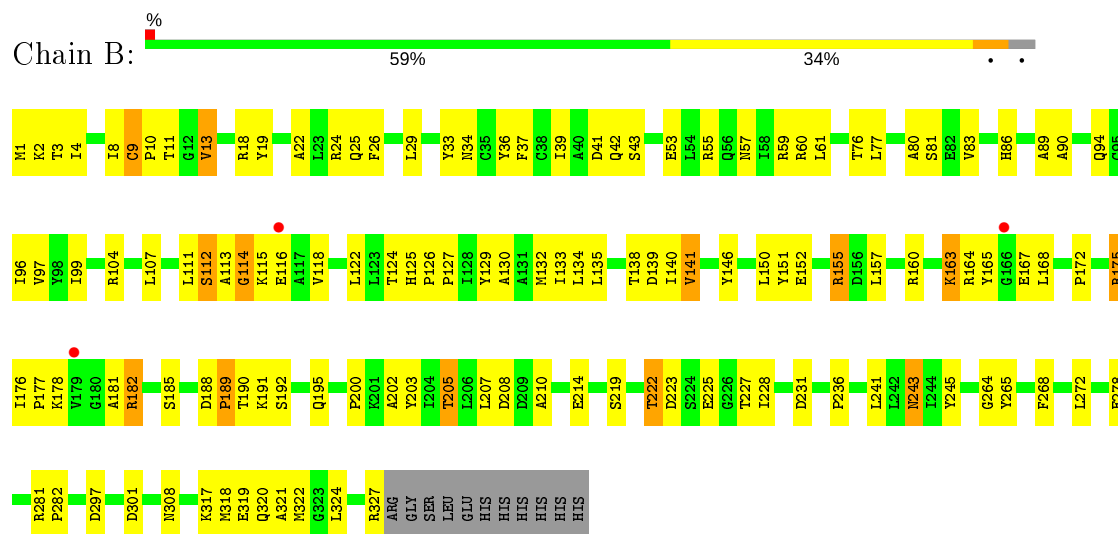
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

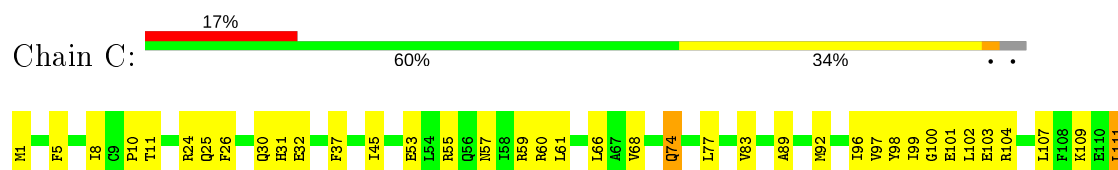
- Molecule 1: De Novo design cysteine esterase FR29



- Molecule 1: De Novo design cysteine esterase FR29



- Molecule 1: De Novo design cysteine esterase FR29



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.75Å 100.76Å 188.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.82 – 2.80 42.82 – 2.80	Depositor EDS
% Data completeness (in resolution range)	90.7 (42.82-2.80) 90.8 (42.82-2.80)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.7.1 _743	Depositor
R, R_{free}	0.214 , 0.290 0.206 , 0.281	Depositor DCC
R_{free} test set	2117 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	39.8	Xtriage
Anisotropy	0.729	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 57.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9968	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.62 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.6741e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.43	0/2638	0.59	0/3554
1	B	0.45	0/2638	0.59	0/3554
1	C	0.38	0/2638	0.55	0/3554
1	D	0.43	0/2177	0.60	0/2933
All	All	0.42	0/10091	0.58	0/13595

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2597	0	2620	69	0
1	B	2597	0	2620	92	0
1	C	2597	0	2620	90	0
1	D	2145	0	2171	87	0
2	A	9	0	0	0	0
2	B	12	0	0	2	0
2	C	6	0	0	1	0
2	D	5	0	0	0	0
All	All	9968	0	10031	318	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 16.

All (318) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:205:THR:HG22	1:D:207:LEU:H	1.12	1.12
1:B:188:ASP:OD1	1:B:190:THR:HG22	1.61	1.01
1:D:213:ILE:HG21	1:D:277:ILE:HG23	1.39	0.98
1:B:175:ARG:HH11	1:B:175:ARG:HG2	1.33	0.93
1:C:205:THR:HG22	1:C:207:LEU:H	1.39	0.88
1:D:276:VAL:HB	1:D:277:ILE:HD12	1.57	0.86
1:C:162:ASN:HA	1:C:166:GLY:O	1.75	0.85
1:D:277:ILE:H	1:D:277:ILE:HD12	1.39	0.84
1:A:205:THR:HG22	1:A:207:LEU:H	1.44	0.81
1:C:215:LYS:O	1:C:219:SER:OG	2.01	0.77
1:C:129:TYR:CZ	1:C:133:ILE:HD11	2.19	0.77
1:D:96:ILE:HG23	1:D:160:ARG:HD2	1.65	0.77
1:D:277:ILE:N	1:D:277:ILE:HD12	2.00	0.77
1:B:222:THR:OG1	1:B:223:ASP:N	2.18	0.76
1:D:176:ILE:HG13	1:D:177:PRO:HD2	1.66	0.76
1:B:155:ARG:NH1	1:B:172:PRO:O	2.17	0.75
1:A:107:LEU:O	1:A:111:LEU:HD23	1.87	0.75
1:B:205:THR:HG22	1:B:207:LEU:H	1.52	0.74
1:C:60:ARG:HG2	1:C:291:MSE:HE2	1.68	0.74
1:A:205:THR:HB	1:A:208:ASP:OD1	1.88	0.73
1:B:43:SER:HB3	1:B:125:HIS:NE2	2.06	0.71
1:C:107:LEU:O	1:C:111:LEU:HD22	1.91	0.71
1:C:60:ARG:HG2	1:C:291:MSE:CE	2.21	0.70
1:C:225:GLU:HG2	1:C:226:GLY:H	1.55	0.69
1:D:83:VAL:HG13	1:D:308:ASN:OD1	1.94	0.68
1:B:324:LEU:HD21	1:D:45:ILE:HD13	1.76	0.68
1:B:86:HIS:CE1	1:B:132:MSE:HG2	2.29	0.67
1:D:14:ILE:HD13	1:D:64:LEU:HG	1.74	0.67
1:D:205:THR:HG22	1:D:207:LEU:N	1.98	0.67
1:C:220:ALA:O	1:C:269:LYS:NZ	2.29	0.66
1:D:66:LEU:CD2	1:D:72:PRO:HG3	2.25	0.66
1:D:96:ILE:CG2	1:D:160:ARG:HD2	2.25	0.66
1:D:115:LYS:HD3	1:D:115:LYS:H	1.61	0.66
1:B:1:MSE:HE2	1:B:33:TYR:C	2.16	0.66
1:D:162:ASN:HA	1:D:166:GLY:O	1.96	0.66
1:A:126:PRO:HB2	1:A:127:PRO:HD3	1.77	0.66
1:D:184:MSE:HG2	1:D:185:SER:H	1.60	0.65
1:B:13:VAL:HG23	1:B:195:GLN:HB2	1.79	0.65
1:A:98:TYR:HD1	1:A:101:GLU:OE2	1.79	0.64
1:A:225:GLU:HG2	1:A:226:GLY:H	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:THR:CG2	1:B:207:LEU:H	2.10	0.64
1:A:116:GLU:CD	1:A:116:GLU:H	2.00	0.64
1:A:133:ILE:HD13	1:A:141:VAL:HG21	1.80	0.63
1:D:203:TYR:C	1:D:204:ILE:HD12	2.18	0.63
1:D:276:VAL:HB	1:D:277:ILE:CD1	2.28	0.63
1:A:175:ARG:O	1:A:177:PRO:HD3	1.98	0.63
1:B:190:THR:HG23	1:B:191:LYS:HG3	1.80	0.63
1:B:115:LYS:HG3	1:B:116:GLU:H	1.63	0.63
1:B:10:PRO:HG3	1:B:41:ASP:HA	1.79	0.63
1:C:104:ARG:HA	1:C:109:LYS:HE3	1.80	0.62
1:A:26:PHE:HA	1:A:29:LEU:HB2	1.81	0.62
1:D:89:ALA:HA	1:D:92:MSE:HE3	1.80	0.62
1:B:26:PHE:HA	1:B:29:LEU:HB2	1.82	0.62
1:A:163:LYS:HD2	1:A:163:LYS:C	2.20	0.61
1:B:55:ARG:NH2	1:D:324:LEU:O	2.33	0.61
1:C:126:PRO:HB2	1:C:127:PRO:HD3	1.83	0.60
1:A:104:ARG:NH2	1:B:113:ALA:H	1.99	0.60
1:C:97:VAL:HG21	1:C:102:LEU:HD21	1.83	0.60
1:B:115:LYS:HD3	1:C:116:GLU:HB2	1.82	0.60
1:D:108:PHE:O	1:D:112:SER:HB3	2.01	0.60
1:D:277:ILE:CD1	1:D:277:ILE:H	2.13	0.60
1:B:133:ILE:HD13	1:B:141:VAL:HG21	1.84	0.60
1:C:205:THR:HG22	1:C:207:LEU:N	2.12	0.60
1:B:9:CYS:HB3	1:B:11:THR:HG22	1.83	0.60
1:B:94:GLN:OE1	1:B:124:THR:HB	2.02	0.59
1:A:184:MSE:HE3	1:A:191:LYS:C	2.23	0.59
1:B:126:PRO:HB2	1:B:127:PRO:HD3	1.82	0.59
1:A:108:PHE:CZ	1:A:123:LEU:HD13	2.38	0.59
1:B:129:TYR:CZ	1:B:133:ILE:HD11	2.37	0.59
1:B:39:ILE:HG12	1:B:61:LEU:HD23	1.83	0.59
1:D:171:ILE:HD12	1:D:171:ILE:N	2.18	0.59
1:D:239:SER:O	1:D:241:LEU:N	2.36	0.59
1:B:11:THR:O	1:B:11:THR:HG23	2.03	0.59
1:D:64:LEU:HD11	1:D:206:LEU:HD12	1.84	0.59
1:A:92:MSE:HE1	1:A:319:GLU:HG3	1.84	0.59
1:B:99:ILE:HG13	1:D:118:VAL:O	2.04	0.57
1:D:246:SER:O	1:D:249:SER:OG	2.22	0.57
1:A:324:LEU:O	1:C:55:ARG:NH2	2.38	0.57
1:D:16:ILE:HD12	1:D:17:GLY:H	1.68	0.57
1:D:197:ASP:HB3	1:D:203:TYR:CE2	2.40	0.57
1:D:25:GLN:HE21	1:D:178:LYS:HD3	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:ALA:HA	1:C:92:MSE:HE3	1.85	0.57
1:A:162:ASN:HA	1:A:166:GLY:O	2.04	0.56
1:B:80:ALA:HB1	2:B:516:HOH:O	2.04	0.56
1:D:183:ILE:HG23	1:D:240:ASN:HD21	1.70	0.56
1:A:104:ARG:HH21	1:B:113:ALA:H	1.53	0.56
1:B:241:LEU:HD22	1:B:272:LEU:HD23	1.88	0.56
1:B:4:ILE:HG12	1:B:140:ILE:HB	1.86	0.56
1:B:175:ARG:NH1	1:B:175:ARG:HG2	2.09	0.56
1:A:222:THR:OG1	1:A:223:ASP:N	2.38	0.56
1:D:13:VAL:HG11	1:D:205:THR:OG1	2.06	0.56
1:C:77:LEU:O	1:C:303:GLY:HA3	2.05	0.55
1:D:276:VAL:CB	1:D:277:ILE:HD12	2.34	0.55
1:A:183:ILE:N	1:A:183:ILE:HD13	2.21	0.55
1:A:301:ASP:OD1	1:C:326:ARG:NH2	2.39	0.55
1:B:115:LYS:HD3	1:C:116:GLU:CB	2.36	0.55
1:C:284:GLN:HA	1:C:287:TYR:HB3	1.88	0.55
1:B:112:SER:O	1:B:113:ALA:HB3	2.06	0.54
1:C:282:PRO:HB2	1:C:286:ARG:NH1	2.21	0.54
1:C:155:ARG:HB2	1:C:155:ARG:NH1	2.22	0.54
1:C:100:GLY:O	1:C:104:ARG:HG3	2.07	0.54
1:B:25:GLN:NE2	1:B:178:LYS:HE2	2.22	0.54
1:B:3:THR:HA	1:B:34:ASN:O	2.08	0.54
1:B:175:ARG:HH11	1:B:175:ARG:CG	2.12	0.54
1:C:97:VAL:HG22	1:C:98:TYR:N	2.23	0.54
1:B:319:GLU:C	1:B:321:ALA:H	2.11	0.54
1:D:245:TYR:O	1:D:245:TYR:HD1	1.90	0.53
1:C:24:ARG:HH22	1:C:247:THR:HB	1.72	0.53
1:C:245:TYR:CD2	1:C:272:LEU:HD13	2.43	0.53
1:C:83:VAL:HG22	1:C:308:ASN:ND2	2.24	0.53
1:C:185:SER:HB3	1:C:188:ASP:O	2.09	0.53
1:C:222:THR:OG1	1:C:223:ASP:N	2.39	0.53
1:D:317:LYS:O	1:D:320:GLN:HB3	2.09	0.53
1:A:129:TYR:CD2	1:A:150:LEU:HG	2.43	0.53
1:D:243:ASN:OD1	1:D:244:ILE:HG13	2.09	0.53
1:D:184:MSE:N	1:D:184:MSE:SE	2.92	0.53
1:A:225:GLU:HG2	1:A:226:GLY:N	2.22	0.53
1:D:13:VAL:HG22	1:D:14:ILE:H	1.74	0.53
1:A:14:ILE:HG13	1:A:64:LEU:HD23	1.90	0.52
1:A:281:ARG:NH1	1:A:282:PRO:HD3	2.24	0.52
1:B:138:THR:HG21	1:B:141:VAL:HG23	1.90	0.52
1:D:100:GLY:O	1:D:104:ARG:HG3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:SER:HB3	1:B:118:VAL:HG21	1.91	0.52
1:B:59:ARG:NH2	1:B:297:ASP:OD2	2.43	0.52
1:A:59:ARG:NH1	1:A:296:LEU:HD23	2.25	0.51
1:B:181:ALA:HB3	1:B:243:ASN:OD1	2.10	0.51
1:D:137:ASN:OD1	1:D:314:MSE:HE3	2.10	0.51
1:A:225:GLU:OE2	1:A:227:THR:OG1	2.29	0.51
1:A:36:TYR:CD1	1:A:76:THR:HB	2.46	0.51
1:C:222:THR:HA	1:C:237:GLY:HA3	1.92	0.51
1:D:199:ASN:HB3	1:D:202:ALA:HB2	1.93	0.51
1:C:103:GLU:HG3	1:C:123:LEU:HD11	1.91	0.51
1:A:99:ILE:HD11	1:C:118:VAL:HB	1.93	0.51
1:D:108:PHE:CZ	1:D:123:LEU:HD13	2.45	0.51
1:B:160:ARG:O	1:B:163:LYS:HD2	2.10	0.51
1:C:26:PHE:O	1:C:30:GLN:HG2	2.11	0.50
1:C:225:GLU:HG2	1:C:226:GLY:N	2.25	0.50
1:A:98:TYR:HB2	1:A:101:GLU:HG3	1.93	0.50
1:D:19:TYR:HA	1:D:23:LEU:HB3	1.92	0.50
1:D:115:LYS:CD	1:D:115:LYS:H	2.24	0.50
1:C:225:GLU:CG	1:C:226:GLY:H	2.23	0.50
1:A:59:ARG:NH2	1:A:297:ASP:OD1	2.45	0.50
1:C:107:LEU:HD13	1:C:107:LEU:O	2.12	0.50
1:C:198:PRO:O	1:C:200:PRO:HD3	2.12	0.50
1:B:175:ARG:NH1	1:B:175:ARG:CG	2.74	0.49
1:B:97:VAL:HG23	1:B:157:LEU:HD21	1.93	0.49
1:B:2:LYS:HE3	1:B:139:ASP:HB3	1.95	0.49
1:B:210:ALA:O	1:B:214:GLU:HG2	2.13	0.49
1:B:319:GLU:HG2	1:B:324:LEU:HB2	1.94	0.49
1:B:42:GLN:NE2	2:B:440:HOH:O	2.35	0.49
1:A:112:SER:OG	1:A:113:ALA:N	2.46	0.49
1:B:150:LEU:O	1:B:150:LEU:HD12	2.11	0.49
1:A:164:ARG:HD2	1:A:165:TYR:CE1	2.48	0.49
1:B:96:ILE:HG22	1:B:160:ARG:HD2	1.94	0.48
1:B:164:ARG:HD2	1:B:165:TYR:CE2	2.48	0.48
1:A:210:ALA:O	1:A:214:GLU:HG2	2.14	0.48
1:C:271:ASP:O	1:C:275:VAL:HG23	2.13	0.48
1:A:125:HIS:O	1:A:128:ILE:HB	2.13	0.48
1:C:129:TYR:OH	1:C:133:ILE:HD11	2.13	0.48
1:C:148:LYS:O	1:C:152:GLU:HB2	2.13	0.48
1:C:53:GLU:HA	1:C:53:GLU:OE1	2.13	0.48
1:A:119:SER:HA	1:C:97:VAL:O	2.13	0.48
1:B:124:THR:HG21	1:D:124:THR:HG21	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:THR:HG22	1:B:207:LEU:N	2.26	0.48
1:C:155:ARG:HH11	1:C:155:ARG:CB	2.27	0.47
1:B:188:ASP:O	1:B:190:THR:N	2.47	0.47
1:C:83:VAL:HA	1:C:308:ASN:HD21	1.80	0.47
1:B:125:HIS:N	1:B:126:PRO:CD	2.78	0.47
1:A:14:ILE:HG21	1:A:206:LEU:HD22	1.97	0.47
1:B:1:MSE:HE2	1:B:33:TYR:O	2.14	0.47
1:D:184:MSE:HG2	1:D:185:SER:N	2.29	0.47
1:D:183:ILE:HG23	1:D:240:ASN:ND2	2.30	0.47
1:B:231:ASP:OD1	1:B:231:ASP:C	2.52	0.47
1:A:94:GLN:OE1	1:A:124:THR:HB	2.14	0.47
1:B:114:GLY:HA2	1:B:115:LYS:HA	1.55	0.47
1:B:86:HIS:ND1	1:B:132:MSE:HG2	2.28	0.47
1:C:184:MSE:HA	1:C:192:SER:HA	1.97	0.47
1:C:205:THR:CG2	1:C:207:LEU:H	2.21	0.47
1:D:15:THR:HG23	1:D:18:ARG:H	1.80	0.47
1:A:41:ASP:OD2	1:A:81:SER:HB3	2.15	0.46
1:C:98:TYR:O	1:C:101:GLU:N	2.46	0.46
1:D:186:LEU:HD11	1:D:204:ILE:CD1	2.44	0.46
1:C:208:ASP:O	1:C:284:GLN:NE2	2.41	0.46
1:D:186:LEU:HB3	1:D:187:VAL:H	1.53	0.46
1:A:14:ILE:HG13	1:A:64:LEU:CD2	2.45	0.46
1:D:105:MSE:O	1:D:106:THR:C	2.54	0.46
1:D:239:SER:N	1:D:241:LEU:HD12	2.31	0.46
1:D:213:ILE:CG2	1:D:277:ILE:HG23	2.27	0.46
1:C:140:ILE:HD13	2:C:466:HOH:O	2.16	0.46
1:D:206:LEU:O	1:D:284:GLN:NE2	2.41	0.45
1:A:164:ARG:O	1:A:166:GLY:N	2.49	0.45
1:B:83:VAL:HA	1:B:308:ASN:HD21	1.80	0.45
1:D:176:ILE:HG13	1:D:177:PRO:CD	2.43	0.45
1:D:204:ILE:HD12	1:D:204:ILE:N	2.31	0.45
1:B:125:HIS:N	1:B:126:PRO:HD2	2.31	0.45
1:C:25:GLN:NE2	1:C:178:LYS:HE2	2.32	0.45
1:C:228:ILE:HG13	1:C:260:TYR:HB3	1.98	0.45
1:B:130:ALA:O	1:B:134:LEU:HG	2.15	0.45
1:B:205:THR:HB	1:B:208:ASP:OD1	2.16	0.45
1:C:136:TYR:CZ	1:C:311:ALA:HB2	2.51	0.45
1:C:155:ARG:HH11	1:C:155:ARG:HB2	1.81	0.45
1:C:221:VAL:O	1:C:237:GLY:HA3	2.17	0.45
1:A:136:TYR:OH	1:A:307:ALA:HB1	2.17	0.45
1:D:56:GLN:O	1:D:59:ARG:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:ALA:O	1:B:94:GLN:HB2	2.17	0.45
1:D:276:VAL:HG12	1:D:277:ILE:HG13	1.99	0.45
1:A:125:HIS:N	1:A:126:PRO:CD	2.80	0.45
1:B:19:TYR:O	1:B:24:ARG:HB2	2.17	0.45
1:C:114:GLY:HA2	1:C:115:LYS:HA	1.70	0.44
1:D:96:ILE:HD12	1:D:96:ILE:HA	1.85	0.44
1:B:176:ILE:HA	1:B:177:PRO:HD2	1.73	0.44
1:B:168:LEU:CD1	1:B:317:LYS:HD3	2.47	0.44
1:C:134:LEU:HB3	1:C:169:PHE:CD1	2.53	0.44
1:C:256:LEU:O	1:C:260:TYR:HB2	2.17	0.44
1:C:66:LEU:HD12	1:C:296:LEU:HD12	1.99	0.44
1:B:281:ARG:N	1:B:282:PRO:HD2	2.32	0.44
1:B:57:ASN:OD1	1:B:60:ARG:NH1	2.50	0.44
1:C:155:ARG:HG2	1:C:156:ASP:N	2.32	0.44
1:A:124:THR:HG21	1:C:124:THR:HG21	1.98	0.44
1:A:124:THR:O	1:A:127:PRO:HD2	2.17	0.44
1:C:282:PRO:HA	1:C:285:GLU:HB2	1.99	0.44
1:D:105:MSE:O	1:D:108:PHE:N	2.45	0.44
1:B:189:PRO:HB2	1:B:236:PRO:HB2	1.98	0.44
1:B:29:LEU:HA	1:B:29:LEU:HD23	1.77	0.44
1:B:168:LEU:HD11	1:B:317:LYS:HD3	1.99	0.44
1:C:248:LEU:HD23	1:C:248:LEU:HA	1.86	0.44
1:B:318:MSE:O	1:B:322:MSE:HG3	2.18	0.44
1:C:168:LEU:HD12	1:C:168:LEU:HA	1.66	0.44
1:D:66:LEU:HD21	1:D:72:PRO:HG3	1.98	0.44
1:D:279:THR:O	1:D:282:PRO:HD2	2.17	0.44
1:A:168:LEU:HD11	1:A:317:LYS:CD	2.47	0.43
1:A:99:ILE:HG13	1:C:118:VAL:O	2.18	0.43
1:B:37:PHE:O	1:B:77:LEU:HD12	2.18	0.43
1:C:140:ILE:HA	1:C:173:GLU:O	2.18	0.43
1:B:319:GLU:C	1:B:321:ALA:N	2.71	0.43
1:C:207:LEU:HD21	1:C:287:TYR:CD2	2.53	0.43
1:D:14:ILE:HD12	1:D:14:ILE:H	1.83	0.43
1:D:79:ILE:HB	1:D:82:GLU:HG3	2.00	0.43
1:C:112:SER:OG	1:C:113:ALA:N	2.48	0.43
1:C:31:HIS:ND1	1:C:74:GLN:OE1	2.49	0.43
1:C:37:PHE:HB2	1:C:77:LEU:CD1	2.49	0.43
1:D:316:ARG:HD2	1:D:327:ARG:NH1	2.33	0.43
1:B:264:GLY:O	1:B:265:TYR:C	2.57	0.43
1:C:227:THR:O	1:C:235:LYS:HE2	2.18	0.43
1:A:160:ARG:HG3	1:A:163:LYS:HZ1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:PRO:HA	1:A:61:LEU:HD22	2.01	0.43
1:D:124:THR:O	1:D:127:PRO:HD2	2.17	0.43
1:A:101:GLU:HG2	1:B:113:ALA:HA	2.00	0.43
1:B:185:SER:OG	1:B:202:ALA:HB1	2.19	0.43
1:D:155:ARG:HB2	1:D:155:ARG:NH1	2.34	0.43
1:A:43:SER:OG	1:A:125:HIS:NE2	2.41	0.43
1:A:278:GLU:HA	1:A:278:GLU:OE1	2.18	0.43
1:A:38:CYS:HB2	1:A:78:PHE:CE1	2.54	0.43
1:B:200:PRO:HA	1:B:203:TYR:CE2	2.54	0.43
1:B:241:LEU:HD22	1:B:272:LEU:CD2	2.49	0.43
1:B:8:ILE:CG2	1:B:61:LEU:HD21	2.49	0.43
1:A:324:LEU:HD21	1:C:45:ILE:HD13	2.00	0.43
1:C:10:PRO:HA	1:C:61:LEU:HD22	2.01	0.43
1:B:43:SER:CB	1:B:125:HIS:NE2	2.79	0.43
1:C:129:TYR:CE2	1:C:133:ILE:CD1	3.02	0.42
1:A:60:ARG:HG2	1:A:287:TYR:OH	2.19	0.42
1:A:96:ILE:HD13	1:A:96:ILE:HA	1.72	0.42
1:C:60:ARG:HG2	1:C:291:MSE:HE1	2.00	0.42
1:D:168:LEU:HD11	1:D:317:LYS:HB3	2.02	0.42
1:A:116:GLU:CD	1:A:116:GLU:N	2.68	0.42
1:D:4:ILE:HG12	1:D:140:ILE:HB	1.99	0.42
1:C:115:LYS:H	1:C:115:LYS:HG3	1.67	0.42
1:C:260:TYR:HA	1:C:263:LYS:HD2	2.01	0.42
1:C:96:ILE:HD13	1:C:96:ILE:HA	1.81	0.42
1:D:18:ARG:O	1:D:22:ALA:HB3	2.20	0.42
1:D:104:ARG:HB2	1:D:104:ARG:HE	1.54	0.42
1:D:186:LEU:HA	1:D:186:LEU:HD23	1.75	0.42
1:A:37:PHE:O	1:A:77:LEU:HD12	2.20	0.42
1:B:36:TYR:CD1	1:B:76:THR:HB	2.55	0.42
1:C:104:ARG:HB2	1:C:104:ARG:HE	1.52	0.42
1:D:239:SER:OG	1:D:240:ASN:N	2.53	0.42
1:A:43:SER:CB	1:A:125:HIS:NE2	2.83	0.41
1:B:182:ARG:HD2	1:B:192:SER:OG	2.19	0.41
1:D:96:ILE:HG23	1:D:160:ARG:CD	2.44	0.41
1:C:274:GLN:O	1:C:278:GLU:HB2	2.20	0.41
1:A:217:ILE:HD13	1:A:217:ILE:HA	1.90	0.41
1:D:130:ALA:O	1:D:134:LEU:HG	2.20	0.41
1:D:281:ARG:N	1:D:282:PRO:HD2	2.35	0.41
1:C:131:ALA:O	1:C:135:LEU:HG	2.21	0.41
1:D:104:ARG:HA	1:D:109:LYS:HE2	2.02	0.41
1:C:37:PHE:HB2	1:C:77:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:284:GLN:HA	1:D:287:TYR:HB3	2.01	0.41
1:B:228:ILE:HD12	1:B:268:PHE:CG	2.56	0.41
1:C:5:PHE:CD1	1:C:5:PHE:C	2.94	0.41
1:B:89:ALA:HB3	1:B:135:LEU:HD11	2.01	0.41
1:B:301:ASP:OD1	1:D:326:ARG:NH2	2.52	0.41
1:C:125:HIS:N	1:C:126:PRO:CD	2.84	0.41
1:C:129:TYR:CE2	1:C:133:ILE:HD11	2.55	0.41
1:C:313:GLU:OE1	1:C:313:GLU:HA	2.20	0.41
1:A:281:ARG:HB3	1:A:282:PRO:HD3	2.03	0.40
1:A:41:ASP:HB3	1:A:58:ILE:HD11	2.02	0.40
1:C:57:ASN:OD1	1:C:60:ARG:NH1	2.54	0.40
1:B:97:VAL:O	1:D:119:SER:HA	2.21	0.40
1:D:125:HIS:N	1:D:126:PRO:CD	2.84	0.40
1:D:201:LYS:HD2	1:D:201:LYS:HA	1.48	0.40
1:D:21:GLY:C	1:D:25:GLN:HE22	2.24	0.40
1:A:225:GLU:OE2	1:A:229:ARG:NH1	2.54	0.40
1:A:320:GLN:O	1:A:320:GLN:HG3	2.21	0.40
1:C:210:ALA:HB1	1:C:277:ILE:HD13	2.02	0.40
1:A:123:LEU:HD22	1:C:99:ILE:HD13	2.03	0.40
1:D:8:ILE:CG2	1:D:61:LEU:HD21	2.51	0.40
1:A:150:LEU:HD12	1:A:150:LEU:HA	1.63	0.40
1:A:56:GLN:OE1	1:A:56:GLN:HA	2.21	0.40
1:D:183:ILE:CG2	1:D:240:ASN:HD21	2.34	0.40
1:A:111:LEU:HD23	1:A:111:LEU:H	1.86	0.40
1:B:41:ASP:OD2	1:B:81:SER:HB3	2.22	0.40
1:B:129:TYR:CZ	1:B:133:ILE:CD1	3.05	0.40
1:B:243:ASN:HA	1:B:243:ASN:HD22	1.62	0.40
1:B:41:ASP:OD1	1:B:42:GLN:N	2.55	0.40
1:C:265:TYR:CD1	1:C:265:TYR:N	2.89	0.40
1:A:55:ARG:NH2	1:C:324:LEU:O	2.47	0.40
1:D:176:ILE:HA	1:D:176:ILE:HD12	1.96	0.40
1:D:36:TYR:CD1	1:D:76:THR:HB	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	325/338 (96%)	299 (92%)	23 (7%)	3 (1%)	17	46
1	B	325/338 (96%)	303 (93%)	18 (6%)	4 (1%)	13	39
1	C	325/338 (96%)	288 (89%)	34 (10%)	3 (1%)	17	46
1	D	261/338 (77%)	233 (89%)	22 (8%)	6 (2%)	6	21
All	All	1236/1352 (91%)	1123 (91%)	97 (8%)	16 (1%)	12	36

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	TYR
1	D	178	LYS
1	D	179	VAL
1	D	240	ASN
1	A	147	GLY
1	C	225	GLU
1	D	208	ASP
1	D	278	GLU
1	C	145	GLU
1	B	189	PRO
1	A	110	GLU
1	B	22	ALA
1	B	320	GLN
1	B	114	GLY
1	C	226	GLY
1	D	45	ILE

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	276/276 (100%)	250 (91%)	26 (9%)	8	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	276/276 (100%)	249 (90%)	27 (10%)	8	24
1	C	276/276 (100%)	257 (93%)	19 (7%)	15	41
1	D	228/276 (83%)	196 (86%)	32 (14%)	3	11
All	All	1056/1104 (96%)	952 (90%)	104 (10%)	8	24

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	9	CYS
1	A	13	VAL
1	A	14	ILE
1	A	53	GLU
1	A	56	GLN
1	A	97	VAL
1	A	116	GLU
1	A	118	VAL
1	A	119	SER
1	A	155	ARG
1	A	163	LYS
1	A	175	ARG
1	A	190	THR
1	A	206	LEU
1	A	219	SER
1	A	223	ASP
1	A	224	SER
1	A	227	THR
1	A	229	ARG
1	A	238	ILE
1	A	245	TYR
1	A	255	GLU
1	A	278	GLU
1	A	298	ARG
1	A	306	LYS
1	B	9	CYS
1	B	13	VAL
1	B	18	ARG
1	B	53	GLU
1	B	104	ARG
1	B	107	LEU
1	B	111	LEU

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Mol	Chain	Res	Type
1	B	112	SER
1	B	122	LEU
1	B	141	VAL
1	B	146	TYR
1	B	151	TYR
1	B	152	GLU
1	B	155	ARG
1	B	163	LYS
1	B	167	GLU
1	B	175	ARG
1	B	182	ARG
1	B	205	THR
1	B	219	SER
1	B	222	THR
1	B	225	GLU
1	B	227	THR
1	B	243	ASN
1	B	245	TYR
1	B	278	GLU
1	B	327	ARG
1	C	1	MSE
1	C	8	ILE
1	C	11	THR
1	C	32	GLU
1	C	59	ARG
1	C	68	VAL
1	C	74	GLN
1	C	111	LEU
1	C	116	GLU
1	C	118	VAL
1	C	119	SER
1	C	128	ILE
1	C	152	GLU
1	C	155	ARG
1	C	182	ARG
1	C	212	THR
1	C	233	GLU
1	C	279	THR
1	C	327	ARG
1	D	6	SER
1	D	8	ILE
1	D	11	THR

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Mol	Chain	Res	Type
1	D	13	VAL
1	D	15	THR
1	D	16	ILE
1	D	55	ARG
1	D	64	LEU
1	D	68	VAL
1	D	96	ILE
1	D	112	SER
1	D	115	LYS
1	D	118	VAL
1	D	146	TYR
1	D	155	ARG
1	D	182	ARG
1	D	183	ILE
1	D	184	MSE
1	D	187	VAL
1	D	199	ASN
1	D	203	TYR
1	D	239	SER
1	D	243	ASN
1	D	245	TYR
1	D	246	SER
1	D	277	ILE
1	D	279	THR
1	D	294	GLU
1	D	302	GLU
1	D	312	SER
1	D	314	MSE
1	D	326	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 [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	317/338 (93%)	-0.12	3 (0%) 84 80	15, 36, 68, 98	0
1	B	317/338 (93%)	-0.20	3 (0%) 84 80	11, 32, 63, 117	0
1	C	317/338 (93%)	0.66	56 (17%) 1 1	14, 54, 141, 162	0
1	D	260/338 (76%)	0.06	15 (5%) 23 15	12, 41, 102, 121	0
All	All	1211/1352 (89%)	0.10	77 (6%) 19 12	11, 38, 114, 162	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	242	LEU	7.0
1	C	218	LYS	7.0
1	C	234	ALA	6.8
1	C	187	VAL	5.1
1	D	245	TYR	4.8
1	C	229	ARG	4.7
1	A	179	VAL	4.7
1	C	188	ASP	4.5
1	C	268	PHE	4.5
1	C	265	TYR	4.5
1	D	213	ILE	4.4
1	C	267	VAL	4.3
1	C	258	ARG	4.0
1	C	222	THR	4.0
1	C	200	PRO	4.0
1	C	264	GLY	4.0
1	C	191	LYS	3.9
1	C	217	ILE	3.9
1	C	238	ILE	3.9
1	C	213	ILE	3.8
1	C	186	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	244	ILE	3.7
1	C	211	LYS	3.7
1	C	221	VAL	3.6
1	C	227	THR	3.5
1	C	203	TYR	3.5
1	C	272	LEU	3.5
1	C	225	GLU	3.3
1	C	226	GLY	3.2
1	D	243	ASN	3.2
1	C	113	ALA	3.2
1	D	242	LEU	3.2
1	C	253	ILE	3.2
1	D	178	LYS	3.2
1	C	212	THR	3.1
1	C	245	TYR	3.1
1	C	189	PRO	3.1
1	C	198	PRO	3.0
1	C	209	ASP	3.0
1	C	233	GLU	3.0
1	A	116	GLU	2.9
1	C	216	LYS	2.9
1	C	256	LEU	2.9
1	D	211	LYS	2.9
1	C	228	ILE	2.9
1	C	208	ASP	2.8
1	C	269	LYS	2.8
1	C	231	ASP	2.8
1	C	214	GLU	2.8
1	D	185	SER	2.8
1	C	281	ARG	2.7
1	C	263	LYS	2.7
1	C	261	GLU	2.7
1	C	199	ASN	2.7
1	D	246	SER	2.7
1	D	180	GLY	2.7
1	B	179	VAL	2.6
1	C	254	GLU	2.6
1	C	196	SER	2.6
1	B	166	GLY	2.4
1	C	202	ALA	2.4
1	D	179	VAL	2.4
1	C	194	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	204	ILE	2.4
1	D	183	ILE	2.4
1	C	219	SER	2.3
1	C	241	LEU	2.3
1	C	274	GLN	2.3
1	D	240	ASN	2.2
1	C	223	ASP	2.2
1	B	116	GLU	2.2
1	A	298	ARG	2.2
1	D	116	GLU	2.1
1	C	197	ASP	2.1
1	C	240	ASN	2.1
1	C	210	ALA	2.1
1	D	248	LEU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.