



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:34 AM GMT

PDB ID : 2R7D  
Title : Crystal structure of ribonuclease II family protein from *Deinococcus radiodurans*, triclinic crystal form. NorthEast Structural Genomics target DrR63  
Authors : Seetharaman, J.; Neely, H.; Forouhar, F.; Wang, D.; Fang, Y.; Cunningham, K.; Ma, L-C.; Xia, R.; Liu, J.; Baran, M.C.; Acton, T.B.; Rost, B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2007-09-07  
Resolution : 1.80 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](mailto: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.

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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

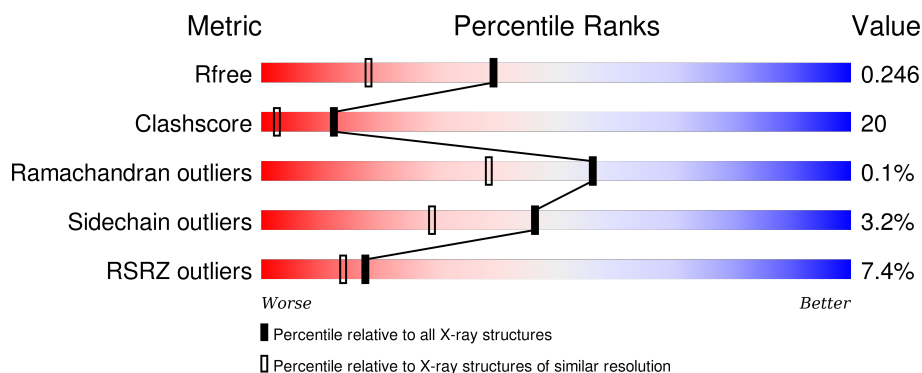
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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  $\geq 3$ , 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	469	<div> <div>3%</div> <div>71%</div> <div>26%</div> <div>..</div> </div>
1	B	469	<div> <div>12%</div> <div>64%</div> <div>32%</div> <div>..</div> </div>
1	C	469	<div> <div>7%</div> <div>70%</div> <div>26%</div> <div>..</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11897 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 Ribonuclease II family protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	Se	0	0	0
			3562	2228	652	671	2	9			
1	B	459	Total	C	N	O	S	Se	0	0	0
			3545	2217	650	667	2	9			
1	C	459	Total	C	N	O	S	Se	0	0	0
			3545	2217	650	667	2	9			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	462	LEU	-	EXPRESSION TAG	UNP Q9RYD0
A	463	GLU	-	EXPRESSION TAG	UNP Q9RYD0
A	464	HIS	-	EXPRESSION TAG	UNP Q9RYD0
A	465	HIS	-	EXPRESSION TAG	UNP Q9RYD0
A	466	HIS	-	EXPRESSION TAG	UNP Q9RYD0
A	467	HIS	-	EXPRESSION TAG	UNP Q9RYD0
A	468	HIS	-	EXPRESSION TAG	UNP Q9RYD0
A	469	HIS	-	EXPRESSION TAG	UNP Q9RYD0
B	462	LEU	-	EXPRESSION TAG	UNP Q9RYD0
B	463	GLU	-	EXPRESSION TAG	UNP Q9RYD0
B	464	HIS	-	EXPRESSION TAG	UNP Q9RYD0
B	465	HIS	-	EXPRESSION TAG	UNP Q9RYD0
B	466	HIS	-	EXPRESSION TAG	UNP Q9RYD0
B	467	HIS	-	EXPRESSION TAG	UNP Q9RYD0
B	468	HIS	-	EXPRESSION TAG	UNP Q9RYD0
B	469	HIS	-	EXPRESSION TAG	UNP Q9RYD0
C	462	LEU	-	EXPRESSION TAG	UNP Q9RYD0
C	463	GLU	-	EXPRESSION TAG	UNP Q9RYD0
C	464	HIS	-	EXPRESSION TAG	UNP Q9RYD0
C	465	HIS	-	EXPRESSION TAG	UNP Q9RYD0
C	466	HIS	-	EXPRESSION TAG	UNP Q9RYD0
C	467	HIS	-	EXPRESSION TAG	UNP Q9RYD0
C	468	HIS	-	EXPRESSION TAG	UNP Q9RYD0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	469	HIS	-	EXPRESSION TAG	UNP Q9RYD0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0

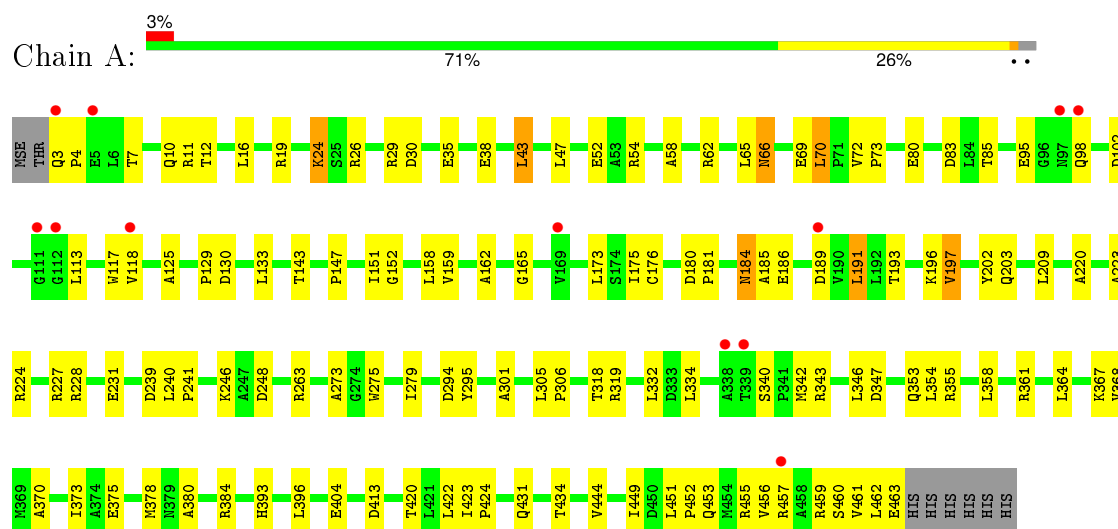
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	536	Total O 536 536	0	0
3	B	303	Total O 303 303	0	0
3	C	403	Total O 403 403	0	0

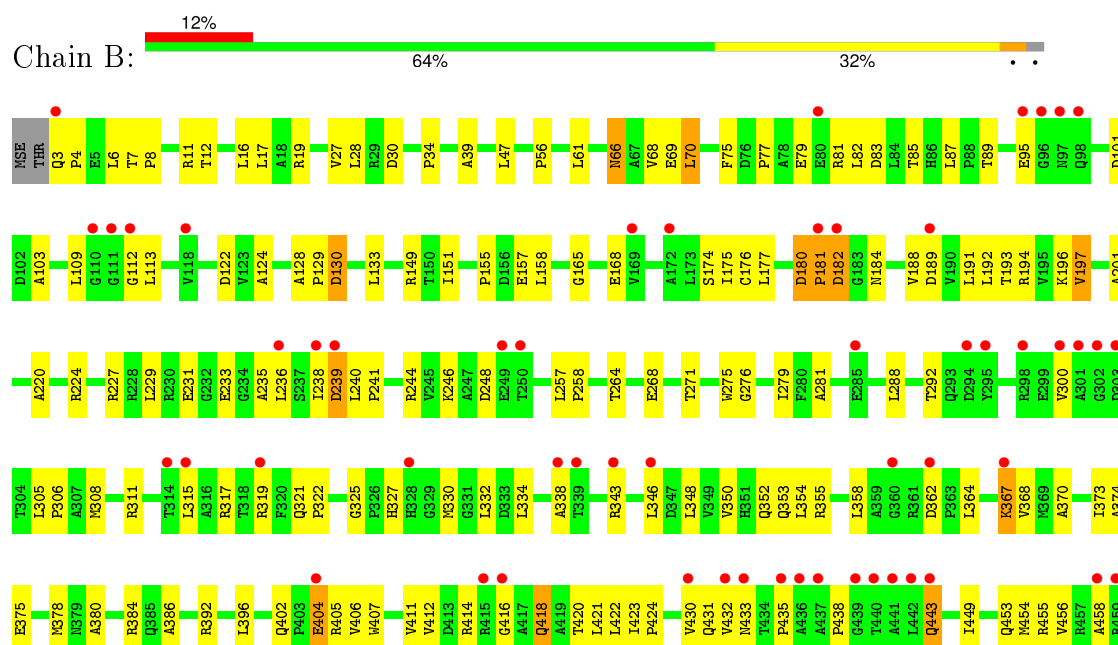
### 3 Residue-property plots

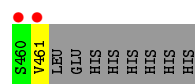
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $RSRZ > 2$ ). 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: Ribonuclease II family protein

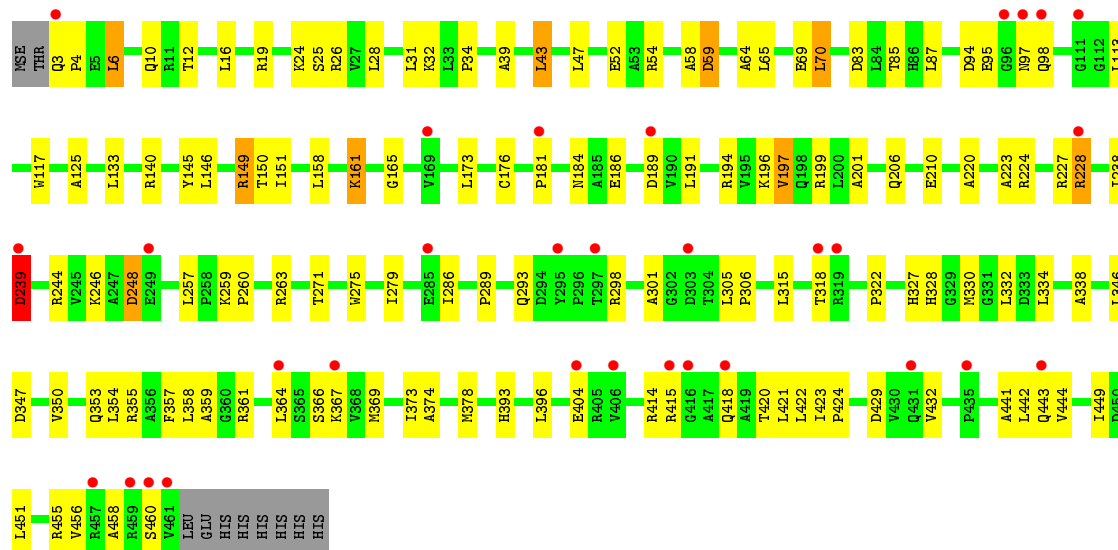


#### • Molecule 1: Ribonuclease II family protein





• Molecule 1: Ribonuclease II family protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.76Å 92.20Å 92.34Å 60.08° 89.57° 71.11°	Depositor
Resolution (Å)	36.68 – 1.80 36.68 – 1.80	Depositor EDS
% Data completeness (in resolution range)	80.7 (36.68-1.80) 76.7 (36.68-1.80)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 1.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.221 , 0.243 0.227 , 0.246	Depositor DCC
$R_{free}$ test set	2691 reflections (1.95%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.4	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 55.5	EDS
Estimated twinning fraction	0.013 for h,h-k,-l 0.016 for -h,-k+l,l 0.008 for -h,-h+k-l,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 308057 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11897	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.33	0/3624	0.62	3/4927 (0.1%)
1	B	0.27	0/3607	0.58	6/4904 (0.1%)
1	C	0.30	0/3607	0.59	2/4904 (0.0%)
All	All	0.30	0/10838	0.60	11/14735 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	239	ASP	CB-CG-OD2	5.27	123.04	118.30
1	B	130	ASP	CB-CG-OD2	5.27	123.04	118.30
1	B	180	ASP	CB-CG-OD2	5.27	123.04	118.30
1	B	362	ASP	CB-CG-OD2	5.25	123.02	118.30
1	B	182	ASP	CB-CG-OD2	5.24	123.01	118.30
1	C	239	ASP	CB-CG-OD2	5.23	123.01	118.30
1	C	248	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	239	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	248	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	248	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	173	LEU	N-CA-C	-5.05	97.35	111.00

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	3562	0	3564	134	0
1	B	3545	0	3547	159	0
1	C	3545	0	3547	127	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	536	0	0	34	0
3	B	303	0	0	24	0
3	C	403	0	0	26	0
All	All	11897	0	10658	420	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 20.

All (420) 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:328:HIS:HB3	3:C:3394:HOH:O	1.60	1.01
1:B:180:ASP:HB2	1:B:181:PRO:HD2	1.41	0.98
1:A:353:GLN:HE22	1:A:364:LEU:H	1.00	0.96
1:B:239:ASP:O	1:B:308:MSE:SE	2.39	0.91
1:A:375:GLU:HA	1:A:378:MSE:HE2	1.53	0.90
1:A:463:GLU:HG2	3:A:1122:HOH:O	1.70	0.90
1:B:321:GLN:HA	3:B:2261:HOH:O	1.72	0.89
1:A:83:ASP:OD2	1:A:85:THR:HG22	1.73	0.89
1:C:52:GLU:HG2	1:C:451:LEU:HD12	1.56	0.88
1:A:370:ALA:HA	3:A:1315:HOH:O	1.74	0.88
1:A:370:ALA:HB3	3:A:1531:HOH:O	1.76	0.85
1:A:26:ARG:HH11	1:A:29:ARG:HH11	1.24	0.85
1:A:176:CYS:SG	1:A:189:ASP:HB2	2.16	0.85
1:C:12:THR:O	1:C:16:LEU:HD23	1.77	0.82
1:B:375:GLU:HA	1:B:378:MSE:HE3	1.61	0.82
1:C:83:ASP:OD2	1:C:85:THR:HG22	1.80	0.82
1:B:124:ALA:HB3	1:B:355:ARG:HD3	1.61	0.81
1:A:159:VAL:HA	1:A:342:MSE:HE1	1.61	0.81
1:C:176:CYS:SG	1:C:189:ASP:HB2	2.22	0.79
1:B:69:GLU:C	1:B:70:LEU:HD12	2.04	0.78
1:A:380:ALA:HB1	1:A:384:ARG:HH12	1.47	0.78
1:B:373:ILE:HB	3:B:2244:HOH:O	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:414:ARG:HB3	1:C:414:ARG:NH1	1.99	0.77
1:B:238:ILE:HG21	1:B:315:LEU:HD11	1.66	0.77
1:B:133:LEU:H	1:B:133:LEU:HD12	1.49	0.76
1:A:380:ALA:HB1	1:A:384:ARG:NH1	2.01	0.76
1:B:353:GLN:HE22	1:B:364:LEU:H	1.31	0.76
1:A:420:THR:HG22	1:A:431:GLN:OE1	1.85	0.76
1:A:162:ALA:HB3	1:A:342:MSE:HE3	1.68	0.76
1:C:279:ILE:HD13	1:C:334:LEU:HD11	1.66	0.76
1:B:70:LEU:CD1	1:B:70:LEU:N	2.50	0.75
1:B:418:GLN:HA	1:B:433:ASN:HA	1.68	0.75
1:B:12:THR:O	1:B:16:LEU:HD23	1.87	0.75
1:A:373:ILE:HB	3:A:1315:HOH:O	1.85	0.75
1:A:367:LYS:HA	3:A:1531:HOH:O	1.85	0.74
1:B:370:ALA:HA	3:B:2244:HOH:O	1.87	0.74
1:A:196:LYS:HG3	1:A:196:LYS:O	1.86	0.74
1:C:414:ARG:HB3	1:C:414:ARG:HH11	1.51	0.74
1:C:196:LYS:HE2	3:C:3032:HOH:O	1.87	0.73
1:C:125:ALA:HB2	1:C:355:ARG:HH12	1.53	0.72
1:C:354:LEU:O	1:C:358:LEU:HD23	1.89	0.72
1:C:228:ARG:NH1	1:C:228:ARG:HB2	2.04	0.72
1:B:176:CYS:SG	1:B:189:ASP:HB2	2.29	0.72
1:A:159:VAL:HA	1:A:342:MSE:CE	2.19	0.72
1:C:196:LYS:O	1:C:196:LYS:HG3	1.89	0.71
1:C:133:LEU:HD12	1:C:133:LEU:H	1.54	0.71
1:A:24:LYS:HB2	1:A:24:LYS:NZ	2.06	0.70
1:B:70:LEU:N	1:B:70:LEU:HD12	2.07	0.70
1:A:118:VAL:CG1	1:A:175:ILE:HB	2.21	0.69
1:A:162:ALA:HB3	1:A:342:MSE:CE	2.22	0.69
1:C:353:GLN:HE22	1:C:364:LEU:H	1.41	0.69
1:C:443:GLN:HA	1:C:443:GLN:HE21	1.58	0.69
1:B:220:ALA:O	1:B:224:ARG:HG3	1.94	0.68
1:A:12:THR:O	1:A:16:LEU:HD23	1.94	0.67
1:B:455:ARG:HD2	3:B:2253:HOH:O	1.95	0.67
1:A:223:ALA:HB1	1:A:227:ARG:HH12	1.60	0.67
1:C:69:GLU:C	1:C:70:LEU:HD12	2.15	0.67
1:A:26:ARG:HH11	1:A:29:ARG:NH1	1.94	0.66
1:C:70:LEU:N	1:C:70:LEU:CD1	2.59	0.66
1:B:129:PRO:HG2	3:B:2119:HOH:O	1.96	0.66
1:B:414:ARG:NH2	1:B:435:PRO:HA	2.11	0.66
1:A:184:ASN:HD22	1:A:185:ALA:N	1.93	0.65
1:B:305:LEU:HB2	1:B:306:PRO:HD3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:ILE:HG13	1:A:456:VAL:HG12	1.77	0.65
1:B:404:GLU:H	1:B:404:GLU:CD	2.00	0.65
1:A:24:LYS:HD3	3:A:1151:HOH:O	1.96	0.65
1:C:369:MSE:O	1:C:373:ILE:HD13	1.96	0.65
1:A:203:GLN:HG3	3:A:1466:HOH:O	1.96	0.65
1:B:95:GLU:HG2	1:B:201:ALA:HB1	1.78	0.65
1:B:165:GLY:HA2	1:B:197:VAL:HG13	1.79	0.65
1:A:364:LEU:HB3	1:A:368:VAL:CG2	2.26	0.65
1:B:244:ARG:HH21	1:B:246:LYS:HZ3	1.44	0.65
1:B:133:LEU:HD12	1:B:133:LEU:N	2.11	0.64
1:A:125:ALA:HB2	1:A:355:ARG:HH12	1.61	0.64
1:A:66:ASN:H	1:A:66:ASN:HD22	1.44	0.64
1:A:353:GLN:HE22	1:A:364:LEU:N	1.83	0.64
1:B:227:ARG:HH22	1:B:268:GLU:HB2	1.63	0.64
1:B:364:LEU:HB3	1:B:368:VAL:CG2	2.28	0.64
1:A:457:ARG:HG3	3:A:1445:HOH:O	1.98	0.64
1:B:124:ALA:CB	1:B:355:ARG:HD3	2.28	0.64
1:C:322:PRO:HD3	3:C:3163:HOH:O	1.96	0.63
1:B:191:LEU:HD23	1:B:191:LEU:N	2.13	0.63
1:C:444:VAL:HG11	1:C:458:ALA:HB1	1.81	0.63
1:C:441:ALA:C	1:C:442:LEU:HD12	2.18	0.63
1:C:449:ILE:HG13	1:C:456:VAL:HG12	1.79	0.63
1:A:180:ASP:OD2	1:A:184:ASN:HB3	1.97	0.63
1:B:240:LEU:HB3	1:B:241:PRO:HD2	1.80	0.62
1:A:69:GLU:C	1:A:70:LEU:HD12	2.19	0.62
1:A:165:GLY:HA2	1:A:197:VAL:HG13	1.82	0.62
1:C:133:LEU:HD12	1:C:133:LEU:N	2.15	0.62
1:A:364:LEU:HB3	1:A:368:VAL:HG21	1.82	0.61
1:A:129:PRO:HG3	3:A:1342:HOH:O	2.01	0.61
1:B:364:LEU:HB3	1:B:368:VAL:HG21	1.81	0.61
1:B:129:PRO:O	1:B:130:ASP:HB3	2.01	0.61
1:A:11:ARG:NH2	3:A:1408:HOH:O	2.33	0.61
1:C:87:LEU:HB3	3:C:3349:HOH:O	1.99	0.61
1:B:30:ASP:HB2	3:B:2088:HOH:O	2.00	0.61
1:B:66:ASN:H	1:B:66:ASN:HD22	1.47	0.61
1:B:85:THR:O	1:B:196:LYS:HE3	2.01	0.61
1:B:422:LEU:O	1:B:424:PRO:HD3	2.01	0.61
1:A:118:VAL:HG11	1:A:273:ALA:CB	2.30	0.61
1:B:4:PRO:HG2	1:B:47:LEU:HD21	1.83	0.61
1:B:244:ARG:HH21	1:B:246:LYS:NZ	1.97	0.60
1:A:54:ARG:HG2	3:A:1341:HOH:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:301:ALA:HB1	3:C:3350:HOH:O	2.01	0.60
1:B:375:GLU:HA	1:B:378:MSE:CE	2.30	0.60
1:A:26:ARG:NH1	1:A:29:ARG:HH11	1.95	0.60
1:B:430:VAL:HG21	1:B:458:ALA:HB2	1.84	0.60
1:C:3:GLN:HB2	1:C:4:PRO:HD3	1.84	0.60
1:B:151:ILE:N	1:B:151:ILE:HD12	2.17	0.59
1:A:305:LEU:HB2	1:A:306:PRO:HD3	1.84	0.59
1:B:227:ARG:NH1	3:B:2072:HOH:O	2.36	0.59
1:A:102:ASP:HB2	3:A:1211:HOH:O	2.02	0.59
1:C:396:LEU:HD22	1:C:456:VAL:CG1	2.33	0.59
1:A:118:VAL:HG11	1:A:273:ALA:HB2	1.85	0.58
1:A:396:LEU:HD22	1:A:456:VAL:CG1	2.33	0.58
1:C:369:MSE:HG3	3:C:3163:HOH:O	2.02	0.58
1:B:330:MSE:HB3	1:B:332:LEU:HD23	1.83	0.58
1:B:319:ARG:HB3	1:B:319:ARG:NH1	2.18	0.58
1:A:318:THR:HG23	3:A:1087:HOH:O	2.02	0.58
1:C:220:ALA:O	1:C:224:ARG:HG3	2.03	0.58
1:B:354:LEU:O	1:B:358:LEU:HD13	2.03	0.58
1:A:24:LYS:HB2	1:A:24:LYS:HZ2	1.68	0.57
1:B:227:ARG:NH2	1:B:264:THR:HG22	2.18	0.57
1:B:8:PRO:HG2	3:B:2047:HOH:O	2.04	0.57
1:A:354:LEU:O	1:A:358:LEU:HD13	2.03	0.57
1:A:420:THR:HG23	3:A:1182:HOH:O	2.02	0.57
1:B:449:ILE:HG13	1:B:456:VAL:HG12	1.84	0.57
1:B:418:GLN:HG3	1:B:431:GLN:HG3	1.87	0.57
1:B:380:ALA:HB1	1:B:384:ARG:NH1	2.19	0.57
1:A:165:GLY:CA	1:A:197:VAL:HG13	2.35	0.57
1:A:453:GLN:HB2	1:A:455:ARG:HH11	1.70	0.57
1:C:24:LYS:HD2	3:C:3329:HOH:O	2.04	0.57
1:A:319:ARG:HD2	3:A:1232:HOH:O	2.04	0.57
1:B:180:ASP:HB2	1:B:181:PRO:CD	2.27	0.56
1:A:184:ASN:HD22	1:A:185:ALA:H	1.52	0.56
1:B:227:ARG:O	1:B:231:GLU:HG3	2.05	0.56
1:A:30:ASP:HB2	3:A:1154:HOH:O	2.05	0.56
1:C:244:ARG:HG3	1:C:257:LEU:HD21	1.86	0.56
1:A:361:ARG:HD3	3:A:1424:HOH:O	2.05	0.56
1:B:402:GLN:OE1	1:B:405:ARG:HD2	2.05	0.56
1:C:443:GLN:HA	1:C:443:GLN:NE2	2.21	0.56
1:C:26:ARG:HH11	1:C:26:ARG:HG2	1.71	0.56
1:A:375:GLU:HA	1:A:378:MSE:CE	2.29	0.56
1:C:396:LEU:HD22	1:C:456:VAL:HG13	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:463:GLU:HB2	3:A:1184:HOH:O	2.05	0.56
1:A:66:ASN:ND2	1:A:66:ASN:H	2.04	0.56
1:C:275:TRP:O	1:C:279:ILE:HG12	2.06	0.56
1:A:70:LEU:N	1:A:70:LEU:CD1	2.68	0.56
1:B:374:ALA:O	1:B:378:MSE:HG3	2.06	0.55
1:A:24:LYS:HG2	1:A:35:GLU:OE1	2.07	0.55
1:A:196:LYS:HE2	3:A:1007:HOH:O	2.06	0.55
1:A:444:VAL:HG23	1:A:459:ARG:O	2.06	0.55
1:C:98:GLN:HA	1:C:199:ARG:HH22	1.71	0.55
1:B:373:ILE:HD12	3:B:2261:HOH:O	2.07	0.55
1:C:54:ARG:HD2	3:C:3145:HOH:O	2.06	0.55
1:B:16:LEU:HD22	1:B:19:ARG:HH21	1.69	0.55
1:A:227:ARG:O	1:A:231:GLU:HG3	2.07	0.55
1:C:3:GLN:HB2	1:C:4:PRO:CD	2.35	0.55
1:A:43:LEU:HD22	1:A:47:LEU:HG	1.89	0.55
1:C:52:GLU:HG2	1:C:451:LEU:CD1	2.35	0.55
1:B:227:ARG:NH1	1:B:268:GLU:OE1	2.40	0.55
1:B:34:PRO:HG2	1:B:39:ALA:HB2	1.88	0.55
1:C:353:GLN:NE2	1:C:364:LEU:HD23	2.21	0.55
1:B:317:ARG:HG2	1:B:317:ARG:HH11	1.71	0.55
1:A:413:ASP:HB3	1:A:420:THR:OG1	2.06	0.55
1:C:70:LEU:HD12	1:C:70:LEU:N	2.22	0.55
1:C:444:VAL:CG1	1:C:458:ALA:HB1	2.37	0.55
1:A:275:TRP:O	1:A:279:ILE:HG12	2.06	0.55
1:B:338:ALA:HB2	1:B:350:VAL:HG11	1.88	0.54
1:C:165:GLY:HA2	1:C:197:VAL:HG13	1.89	0.54
1:C:59:ASP:OD1	1:C:65:LEU:HD11	2.06	0.54
1:A:3:GLN:N	1:A:4:PRO:HD2	2.23	0.54
1:A:340:SER:HB3	1:A:343:ARG:HG2	1.90	0.54
1:B:348:LEU:O	1:B:352:GLN:HG3	2.07	0.54
1:C:228:ARG:HH11	1:C:228:ARG:HB2	1.71	0.54
1:B:432:VAL:HG22	1:B:458:ALA:HB3	1.90	0.54
1:B:11:ARG:NH2	3:B:2226:HOH:O	2.40	0.54
1:B:279:ILE:HD13	1:B:334:LEU:HD11	1.89	0.54
1:B:7:THR:HB	1:B:8:PRO:HD2	1.89	0.54
1:C:327:HIS:CE1	1:C:330:MSE:HG3	2.44	0.53
1:A:19:ARG:NH2	1:A:54:ARG:HH22	2.06	0.53
1:B:449:ILE:HD12	1:B:449:ILE:N	2.23	0.53
1:C:420:THR:HB	3:C:3247:HOH:O	2.09	0.53
1:A:118:VAL:HG13	1:A:118:VAL:O	2.08	0.53
1:C:191:LEU:HD23	1:C:191:LEU:N	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:GLY:CA	1:B:197:VAL:HG13	2.39	0.53
1:B:174:SER:C	1:B:175:ILE:HD12	2.29	0.53
1:C:69:GLU:O	1:C:70:LEU:HD12	2.09	0.53
1:A:69:GLU:O	1:A:70:LEU:HD12	2.09	0.53
1:B:396:LEU:HD22	1:B:456:VAL:CG1	2.39	0.53
1:C:364:LEU:HD22	1:C:364:LEU:H	1.74	0.52
1:A:346:LEU:HB2	3:A:1356:HOH:O	2.09	0.52
1:A:158:LEU:HD23	1:A:342:MSE:CE	2.38	0.52
1:C:133:LEU:H	1:C:133:LEU:CD1	2.20	0.52
1:A:396:LEU:HD22	1:A:456:VAL:HG11	1.91	0.52
1:B:227:ARG:NH1	3:B:2138:HOH:O	2.42	0.52
1:C:449:ILE:HG13	1:C:456:VAL:CG1	2.39	0.52
1:B:334:LEU:N	1:B:334:LEU:HD23	2.25	0.52
1:C:184:ASN:ND2	3:C:3378:HOH:O	2.40	0.52
1:B:418:GLN:N	1:B:418:GLN:HE21	2.07	0.52
1:B:180:ASP:C	1:B:182:ASP:H	2.12	0.52
1:B:180:ASP:OD2	1:B:184:ASN:HB2	2.08	0.52
1:A:453:GLN:HG3	1:A:455:ARG:NH1	2.25	0.51
1:A:118:VAL:HG13	1:A:175:ILE:HB	1.93	0.51
1:A:151:ILE:HD12	1:A:151:ILE:N	2.25	0.51
1:C:140:ARG:NH1	3:C:3325:HOH:O	2.43	0.51
1:C:95:GLU:HG3	3:C:3025:HOH:O	2.09	0.51
1:A:361:ARG:HH11	1:A:361:ARG:HG3	1.76	0.51
1:B:300:VAL:HG23	1:B:311:ARG:HE	1.76	0.51
1:B:133:LEU:CD1	1:B:133:LEU:H	2.18	0.51
1:B:85:THR:C	3:B:2300:HOH:O	2.48	0.51
1:C:194:ARG:HG3	1:C:194:ARG:HH11	1.76	0.51
1:B:89:THR:HG22	1:B:103:ALA:HB1	1.93	0.51
1:B:453:GLN:NE2	3:B:2263:HOH:O	2.43	0.51
1:B:406:VAL:HG12	1:B:443:GLN:HG3	1.93	0.50
1:B:180:ASP:OD1	1:B:182:ASP:HB2	2.11	0.50
1:B:327:HIS:CE1	1:B:330:MSE:HG3	2.46	0.50
1:C:338:ALA:HB2	1:C:350:VAL:HG11	1.92	0.50
1:C:133:LEU:HD11	3:C:3119:HOH:O	2.11	0.50
1:B:129:PRO:O	1:B:130:ASP:CB	2.59	0.50
1:A:355:ARG:NH2	3:A:1020:HOH:O	2.45	0.50
1:A:117:TRP:HZ3	1:A:191:LEU:HD11	1.77	0.50
1:A:147:PRO:O	1:A:246:LYS:HD3	2.10	0.50
1:B:149:ARG:HD2	1:B:151:ILE:HD11	1.94	0.50
1:B:95:GLU:HG3	3:B:2101:HOH:O	2.12	0.50
1:B:405:ARG:HD3	1:B:407:TRP:CZ2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:GLN:N	1:B:4:PRO:HD2	2.27	0.50
1:B:412:VAL:HG22	1:B:420:THR:O	2.12	0.50
1:C:85:THR:O	1:C:196:LYS:HD2	2.11	0.50
1:A:191:LEU:HD13	1:A:193:THR:CG2	2.42	0.50
1:B:392:ARG:NH1	1:B:454:MSE:HB3	2.27	0.50
1:A:52:GLU:O	1:A:393:HIS:HE1	1.95	0.50
1:A:449:ILE:HG13	1:A:456:VAL:CG1	2.42	0.49
1:C:145:TYR:HE1	3:C:3360:HOH:O	1.95	0.49
1:B:109:LEU:HB2	1:B:113:LEU:HB3	1.94	0.49
1:C:238:ILE:HD13	1:C:315:LEU:HD11	1.94	0.49
1:A:133:LEU:HD13	3:A:1019:HOH:O	2.12	0.49
1:C:85:THR:O	1:C:196:LYS:CD	2.61	0.49
1:A:158:LEU:HD23	1:A:342:MSE:HE2	1.94	0.49
1:A:460:SER:HB3	3:A:1184:HOH:O	2.12	0.49
1:B:82:LEU:HB2	3:B:2063:HOH:O	2.11	0.49
1:C:113:LEU:HD11	1:C:186:GLU:OE1	2.13	0.49
1:C:364:LEU:HD22	1:C:364:LEU:N	2.28	0.49
1:B:69:GLU:O	1:B:70:LEU:HD12	2.12	0.48
1:B:8:PRO:HD3	3:B:2226:HOH:O	2.12	0.48
1:A:453:GLN:HG3	1:A:455:ARG:HH12	1.78	0.48
1:C:359:ALA:HB3	1:C:361:ARG:HG3	1.95	0.48
1:A:191:LEU:HD13	1:A:193:THR:HG23	1.95	0.48
1:A:240:LEU:HB3	1:A:241:PRO:HD2	1.96	0.48
1:B:229:LEU:O	1:B:233:GLU:HG2	2.14	0.48
1:B:124:ALA:HB3	1:B:355:ARG:CD	2.38	0.48
1:B:311:ARG:O	1:B:315:LEU:HG	2.12	0.48
1:C:364:LEU:CD2	1:C:364:LEU:H	2.26	0.48
1:B:414:ARG:HG2	1:B:416:GLY:H	1.77	0.48
1:C:165:GLY:CA	1:C:197:VAL:HG13	2.43	0.48
1:C:361:ARG:HG2	1:C:361:ARG:HH11	1.78	0.48
1:B:180:ASP:O	1:B:182:ASP:N	2.46	0.48
1:A:26:ARG:NH1	3:A:1382:HOH:O	2.43	0.48
1:B:175:ILE:HD12	1:B:175:ILE:N	2.29	0.48
1:A:393:HIS:HD2	3:A:1061:HOH:O	1.97	0.48
1:A:58:ALA:CB	1:A:65:LEU:HD21	2.43	0.48
1:C:151:ILE:HD12	1:C:151:ILE:N	2.28	0.48
1:B:317:ARG:NH1	1:B:317:ARG:HG2	2.28	0.48
1:B:418:GLN:HB3	1:B:433:ASN:CG	2.33	0.48
1:B:112:GLY:O	1:B:181:PRO:HD3	2.13	0.48
1:A:355:ARG:CZ	3:A:1020:HOH:O	2.60	0.48
1:B:194:ARG:HH11	1:B:194:ARG:HG3	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:330:MSE:O	1:C:332:LEU:HD13	2.14	0.47
1:C:374:ALA:O	1:C:378:MSE:HG2	2.15	0.47
1:A:364:LEU:HB3	1:A:368:VAL:HG23	1.96	0.47
1:C:414:ARG:HH11	1:C:414:ARG:CB	2.25	0.47
1:B:396:LEU:HD22	1:B:456:VAL:HG11	1.96	0.47
1:C:422:LEU:O	1:C:424:PRO:HD3	2.15	0.47
1:B:367:LYS:HD2	1:B:367:LYS:C	2.34	0.47
1:B:75:PHE:O	1:B:77:PRO:HD3	2.14	0.47
1:B:418:GLN:N	1:B:418:GLN:NE2	2.63	0.47
1:C:125:ALA:CB	1:C:355:ARG:HH12	2.24	0.47
1:A:70:LEU:N	1:A:70:LEU:HD12	2.28	0.47
1:A:7:THR:O	1:A:11:ARG:HG3	2.14	0.47
1:A:461:VAL:HG22	1:A:462:LEU:HG	1.97	0.47
1:C:421:LEU:HD13	1:C:432:VAL:HB	1.97	0.47
1:A:158:LEU:HD23	1:A:158:LEU:C	2.34	0.47
1:C:58:ALA:CB	1:C:65:LEU:HD21	2.45	0.47
1:C:305:LEU:N	1:C:306:PRO:HD2	2.30	0.47
1:C:94:ASP:OD2	1:C:263:ARG:NH2	2.46	0.47
1:B:227:ARG:HG3	1:B:227:ARG:HH11	1.80	0.47
1:B:240:LEU:HB3	1:B:241:PRO:CD	2.45	0.47
1:B:83:ASP:OD2	1:B:85:THR:HG22	2.15	0.47
1:C:415:ARG:O	1:C:418:GLN:HB2	2.15	0.47
1:A:158:LEU:CD2	1:A:342:MSE:HE2	2.45	0.46
1:C:43:LEU:HD22	1:C:47:LEU:HG	1.97	0.46
1:C:271:THR:HG23	1:C:330:MSE:HE2	1.97	0.46
1:B:66:ASN:ND2	1:B:66:ASN:H	2.11	0.46
1:C:181:PRO:HD2	3:C:3339:HOH:O	2.14	0.46
1:C:423:ILE:N	1:C:423:ILE:HD12	2.30	0.46
1:B:411:VAL:CG1	1:B:438:PRO:HA	2.46	0.46
1:B:373:ILE:CD1	3:B:2261:HOH:O	2.63	0.46
1:A:38:GLU:HG2	3:A:1332:HOH:O	2.14	0.46
1:C:393:HIS:HD2	3:C:3125:HOH:O	1.97	0.46
1:C:318:THR:HG23	3:C:3115:HOH:O	2.14	0.46
1:B:17:LEU:HD23	1:B:28:LEU:HD12	1.97	0.46
1:C:34:PRO:HG2	1:C:39:ALA:HB2	1.97	0.46
1:B:292:THR:HG22	1:B:325:GLY:O	2.16	0.46
1:C:228:ARG:HH11	1:C:228:ARG:CB	2.29	0.46
1:B:271:THR:HG23	1:B:330:MSE:HE2	1.97	0.46
1:B:380:ALA:HB1	1:B:384:ARG:HH12	1.80	0.46
1:B:87:LEU:HB3	3:B:2303:HOH:O	2.14	0.46
1:B:275:TRP:O	1:B:279:ILE:HG12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:25:SER:OG	1:C:28:LEU:HD13	2.15	0.45
1:A:98:GLN:O	1:A:98:GLN:HG2	2.16	0.45
1:C:196:LYS:HE3	3:C:3327:HOH:O	2.15	0.45
1:B:70:LEU:N	1:B:70:LEU:HD13	2.32	0.45
1:B:418:GLN:CA	1:B:433:ASN:HA	2.43	0.45
1:C:248:ASP:OD1	1:C:248:ASP:C	2.55	0.45
1:B:192:LEU:HD22	1:B:355:ARG:NH1	2.31	0.45
1:C:449:ILE:HA	1:C:456:VAL:HG12	1.98	0.45
1:B:61:LEU:HD12	1:B:386:ALA:HB2	1.98	0.45
1:B:158:LEU:HD23	1:B:158:LEU:C	2.36	0.45
1:B:343:ARG:NH2	3:B:2287:HOH:O	2.49	0.45
1:B:238:ILE:HD13	1:B:315:LEU:CD1	2.47	0.45
1:A:165:GLY:N	1:A:197:VAL:HG13	2.32	0.45
1:B:330:MSE:O	1:B:332:LEU:HD22	2.16	0.45
1:B:281:ALA:HB2	1:B:288:LEU:HD11	1.99	0.45
1:B:423:ILE:HD12	1:B:423:ILE:N	2.32	0.45
1:A:279:ILE:CD1	1:A:334:LEU:HD21	2.47	0.45
1:B:443:GLN:HB3	1:B:461:VAL:HG12	1.98	0.45
1:A:113:LEU:HD11	1:A:186:GLU:OE2	2.17	0.45
1:A:52:GLU:HG3	3:A:1200:HOH:O	2.17	0.44
1:B:406:VAL:CG1	1:B:443:GLN:HG3	2.46	0.44
1:C:353:GLN:HA	1:C:353:GLN:NE2	2.32	0.44
1:B:128:ALA:HB1	1:B:129:PRO:HD2	1.99	0.44
1:C:455:ARG:NE	3:C:3317:HOH:O	2.50	0.44
1:C:346:LEU:HB2	3:C:3101:HOH:O	2.16	0.44
1:B:188:VAL:HG11	1:B:276:GLY:CA	2.48	0.44
1:B:353:GLN:NE2	1:B:364:LEU:H	2.06	0.44
1:B:180:ASP:CB	1:B:181:PRO:HD2	2.25	0.44
1:A:7:THR:OG1	1:A:10:GLN:HG3	2.18	0.44
1:C:165:GLY:N	1:C:197:VAL:HG13	2.33	0.44
1:A:62:ARG:HH11	1:A:62:ARG:HG3	1.82	0.44
1:A:95:GLU:HG3	3:A:1105:HOH:O	2.17	0.44
1:B:180:ASP:C	1:B:182:ASP:N	2.70	0.44
1:C:150:THR:HG21	3:C:3360:HOH:O	2.18	0.44
1:A:228:ARG:HD3	3:A:1421:HOH:O	2.17	0.44
1:A:453:GLN:HB2	1:A:455:ARG:NH1	2.33	0.43
1:C:64:ALA:HB3	3:C:3325:HOH:O	2.18	0.43
1:B:418:GLN:NE2	1:B:418:GLN:H	2.15	0.43
1:C:367:LYS:HD2	1:C:367:LYS:O	2.18	0.43
1:B:157:GLU:HG3	3:B:2202:HOH:O	2.18	0.43
1:B:300:VAL:CG2	1:B:311:ARG:HE	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:298:ARG:HG3	1:C:298:ARG:HH11	1.83	0.43
1:A:72:VAL:HA	1:A:73:PRO:HD3	1.86	0.43
1:C:322:PRO:HB2	1:C:366:SER:HB2	2.00	0.43
1:B:430:VAL:HG21	1:B:458:ALA:CB	2.49	0.43
1:B:87:LEU:N	3:B:2300:HOH:O	2.52	0.43
1:B:421:LEU:HD12	1:B:421:LEU:N	2.34	0.43
1:C:334:LEU:N	1:C:334:LEU:HD23	2.34	0.43
1:C:223:ALA:HB1	1:C:227:ARG:HH12	1.84	0.43
1:C:26:ARG:HG2	1:C:26:ARG:NH1	2.32	0.43
1:C:429:ASP:HB2	3:C:3247:HOH:O	2.19	0.43
1:B:177:LEU:HD12	1:B:177:LEU:N	2.34	0.43
1:C:173:LEU:HD23	1:C:173:LEU:C	2.39	0.43
1:B:81:ARG:HD3	1:B:193:THR:O	2.19	0.43
1:A:455:ARG:HG2	1:A:455:ARG:HH11	1.83	0.42
1:C:52:GLU:O	1:C:393:HIS:HE1	2.02	0.42
1:C:69:GLU:C	1:C:70:LEU:CD1	2.85	0.42
1:B:79:GLU:OE2	1:B:192:LEU:HD13	2.19	0.42
1:A:373:ILE:HD12	3:A:1315:HOH:O	2.18	0.42
1:A:118:VAL:HG23	3:A:1211:HOH:O	2.17	0.42
1:C:146:LEU:HD12	1:C:149:ARG:NH1	2.35	0.42
1:A:294:ASP:CG	1:A:295:TYR:H	2.23	0.42
1:A:451:LEU:HB2	1:A:452:PRO:HD3	2.01	0.42
1:B:231:GLU:HA	1:B:235:ALA:HB3	2.01	0.42
1:A:165:GLY:N	1:A:197:VAL:CG1	2.83	0.42
1:C:94:ASP:O	1:C:97:ASN:OD1	2.37	0.42
1:B:346:LEU:HD13	3:B:2010:HOH:O	2.18	0.42
1:C:293:GLN:CG	3:C:3394:HOH:O	2.67	0.42
1:C:443:GLN:O	1:C:460:SER:HA	2.19	0.42
1:A:180:ASP:HB2	1:A:181:PRO:CD	2.49	0.42
1:A:422:LEU:O	1:A:424:PRO:HD3	2.20	0.42
1:C:286:ILE:HD11	1:C:357:PHE:CG	2.55	0.42
1:B:319:ARG:HH11	1:B:319:ARG:CB	2.32	0.42
1:C:31:LEU:O	1:C:32:LYS:HB2	2.19	0.42
1:B:322:PRO:HD3	3:B:2261:HOH:O	2.20	0.41
1:B:81:ARG:NH2	1:B:122:ASP:OD2	2.53	0.41
1:C:161:LYS:CB	1:C:161:LYS:NZ	2.82	0.41
1:B:68:VAL:HG12	1:B:155:PRO:HG3	2.01	0.41
1:B:414:ARG:HG3	1:B:414:ARG:NH1	2.34	0.41
1:C:19:ARG:NH1	3:C:3030:HOH:O	2.52	0.41
1:A:423:ILE:N	1:A:423:ILE:HD12	2.35	0.41
1:A:396:LEU:HD22	1:A:456:VAL:HG13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:330:MSE:CB	1:B:332:LEU:HD23	2.47	0.41
1:C:259:LYS:N	1:C:260:PRO:CD	2.83	0.41
1:B:236:LEU:HD12	3:B:2234:HOH:O	2.20	0.41
1:C:289:PRO:HB3	1:C:373:ILE:HD11	2.03	0.41
1:A:294:ASP:CG	1:A:295:TYR:N	2.74	0.41
1:A:184:ASN:ND2	1:A:185:ALA:N	2.65	0.41
1:C:158:LEU:C	1:C:158:LEU:HD23	2.41	0.41
1:B:257:LEU:HA	1:B:258:PRO:HD3	1.95	0.41
1:A:202:TYR:CD1	1:A:263:ARG:HA	2.56	0.41
1:C:449:ILE:N	1:C:449:ILE:HD12	2.35	0.41
1:B:6:LEU:HD21	1:B:27:VAL:HG11	2.03	0.41
1:A:343:ARG:NH1	3:A:1284:HOH:O	2.54	0.41
1:A:117:TRP:CZ3	1:A:191:LEU:HD11	2.56	0.41
1:C:161:LYS:NZ	3:C:3399:HOH:O	2.47	0.41
1:C:206:GLN:O	1:C:210:GLU:HG3	2.21	0.41
1:A:24:LYS:HB2	1:A:24:LYS:HZ3	1.83	0.41
1:A:449:ILE:N	1:A:449:ILE:HD12	2.36	0.41
1:A:449:ILE:HA	1:A:456:VAL:HG12	2.03	0.41
1:A:220:ALA:O	1:A:224:ARG:HG3	2.20	0.41
1:B:191:LEU:CD2	1:B:191:LEU:N	2.81	0.40
1:C:355:ARG:HH11	1:C:355:ARG:HG3	1.87	0.40
1:C:95:GLU:HG2	1:C:201:ALA:HB1	2.02	0.40
1:C:6:LEU:HG	1:C:10:GLN:HB3	2.03	0.40
1:A:143:THR:HG23	1:A:152:GLY:HA2	2.03	0.40
1:A:301:ALA:HB1	3:A:1411:HOH:O	2.19	0.40
1:A:279:ILE:HD13	1:A:334:LEU:HD21	2.03	0.40
1:C:194:ARG:NH1	1:C:194:ARG:HG3	2.36	0.40
1:C:117:TRP:CE3	1:C:176:CYS:HB3	2.57	0.40
1:B:411:VAL:HG12	1:B:438:PRO:HA	2.03	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	459/469 (98%)	450 (98%)	9 (2%)	0	100	100
1	B	457/469 (97%)	441 (96%)	15 (3%)	1 (0%)	52	35
1	C	457/469 (97%)	450 (98%)	6 (1%)	1 (0%)	52	35
All	All	1373/1407 (98%)	1341 (98%)	30 (2%)	2 (0%)	56	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	239	ASP
1	B	181	PRO

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	371/369 (100%)	357 (96%)	14 (4%)	40	22
1	B	369/369 (100%)	359 (97%)	10 (3%)	52	36
1	C	369/369 (100%)	357 (97%)	12 (3%)	45	27
All	All	1109/1107 (100%)	1073 (97%)	36 (3%)	46	29

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LYS
1	A	43	LEU
1	A	66	ASN
1	A	70	LEU
1	A	80	GLU
1	A	130	ASP
1	A	184	ASN
1	A	191	LEU
1	A	197	VAL
1	A	209	LEU
1	A	332	LEU

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Mol	Chain	Res	Type
1	A	347	ASP
1	A	404	GLU
1	A	434	THR
1	B	56	PRO
1	B	66	ASN
1	B	70	LEU
1	B	101	ASP
1	B	168	GLU
1	B	197	VAL
1	B	367	LYS
1	B	404	GLU
1	B	418	GLN
1	B	443	GLN
1	C	6	LEU
1	C	43	LEU
1	C	59	ASP
1	C	70	LEU
1	C	149	ARG
1	C	161	LYS
1	C	197	VAL
1	C	228	ARG
1	C	239	ASP
1	C	246	LYS
1	C	347	ASP
1	C	404	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	97	ASN
1	A	184	ASN
1	A	321	GLN
1	A	353	GLN
1	A	393	HIS
1	A	433	ASN
1	B	66	ASN
1	B	97	ASN
1	B	184	ASN
1	B	353	GLN
1	B	393	HIS
1	B	418	GLN

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Mol	Chain	Res	Type
1	B	431	GLN
1	C	66	ASN
1	C	97	ASN
1	C	98	GLN
1	C	184	ASN
1	C	353	GLN
1	C	393	HIS
1	C	443	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](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	452/469 (96%)	0.17	12 (2%) 58 53	10, 24, 41, 58	0
1	B	450/469 (95%)	0.73	57 (12%) 5 4	15, 37, 59, 66	0
1	C	450/469 (95%)	0.37	31 (6%) 20 16	15, 30, 50, 59	0
All	All	1352/1407 (96%)	0.42	100 (7%) 17 14	10, 30, 53, 66	0

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	461	VAL	6.9
1	C	461	VAL	6.2
1	B	435	PRO	6.0
1	B	3	GLN	5.5
1	A	3	GLN	4.6
1	C	189	ASP	4.3
1	B	416	GLY	4.2
1	B	415	ARG	4.2
1	B	303	ASP	4.1
1	B	440	THR	4.0
1	C	435	PRO	3.8
1	B	295	TYR	3.8
1	B	319	ARG	3.8
1	B	169	VAL	3.7
1	C	181	PRO	3.7
1	A	189	ASP	3.6
1	B	302	GLY	3.5
1	B	443	GLN	3.4
1	C	111	GLY	3.4
1	B	181	PRO	3.3
1	C	364	LEU	3.3
1	B	98	GLN	3.3
1	C	98	GLN	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	459	ARG	3.3
1	B	249	GLU	3.2
1	B	189	ASP	3.2
1	B	315	LEU	3.2
1	A	118	VAL	3.2
1	C	295	TYR	3.1
1	B	285	GLU	3.1
1	B	460	SER	3.1
1	B	97	ASN	3.1
1	B	442	LEU	3.0
1	A	98	GLN	3.0
1	B	300	VAL	3.0
1	B	314	THR	3.0
1	B	96	GLY	3.0
1	B	236	LEU	3.0
1	B	301	ALA	2.9
1	B	250	THR	2.9
1	B	436	ALA	2.9
1	C	249	GLU	2.9
1	B	118	VAL	2.9
1	C	228	ARG	2.8
1	C	459	ARG	2.8
1	A	112	GLY	2.8
1	C	3	GLN	2.8
1	B	439	GLY	2.7
1	A	111	GLY	2.7
1	C	297	THR	2.7
1	C	415	ARG	2.7
1	B	433	ASN	2.6
1	C	303	ASP	2.6
1	B	430	VAL	2.6
1	B	404	GLU	2.6
1	B	238	ILE	2.6
1	C	431	GLN	2.6
1	B	110	GLY	2.6
1	A	339	THR	2.6
1	B	339	THR	2.6
1	C	416	GLY	2.5
1	B	338	ALA	2.5
1	C	443	GLN	2.5
1	A	97	ASN	2.4
1	C	97	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	298	ARG	2.4
1	B	112	GLY	2.4
1	B	343	ARG	2.4
1	B	172	ALA	2.3
1	B	294	ASP	2.3
1	A	169	VAL	2.3
1	B	95	GLU	2.3
1	C	96	GLY	2.3
1	B	362	ASP	2.3
1	B	437	ALA	2.3
1	C	460	SER	2.2
1	B	458	ALA	2.2
1	C	457	ARG	2.2
1	C	404	GLU	2.2
1	A	338	ALA	2.2
1	B	432	VAL	2.2
1	B	360	GLY	2.2
1	B	367	LYS	2.2
1	A	5	GLU	2.1
1	B	346	LEU	2.1
1	C	418	GLN	2.1
1	C	406	VAL	2.1
1	C	367	LYS	2.1
1	C	285	GLU	2.1
1	B	441	ALA	2.1
1	C	169	VAL	2.1
1	B	182	ASP	2.1
1	A	457	ARG	2.1
1	B	111	GLY	2.1
1	B	239	ASP	2.1
1	C	318	THR	2.0
1	B	80	GLU	2.0
1	C	239	ASP	2.0
1	C	319	ARG	2.0
1	B	328	HIS	2.0

## 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 [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	C	3000	1/1	0.98	0.05	-	20,20,20,20	0
2	MG	B	2000	1/1	0.96	0.06	-	19,19,19,19	0
2	MG	A	1000	1/1	0.97	0.15	-	20,20,20,20	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.