



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:49 PM GMT

PDB ID : 3N4D  
Title : Crystal structure of Cg10062 inactivated by(R)-oxirane-2-carboxylate  
Authors : Guo, Y.; Robertson, B.A.; Hackert, M.L.; Whitman, C.P.  
Deposited on : 2010-05-21  
Resolution : 2.54 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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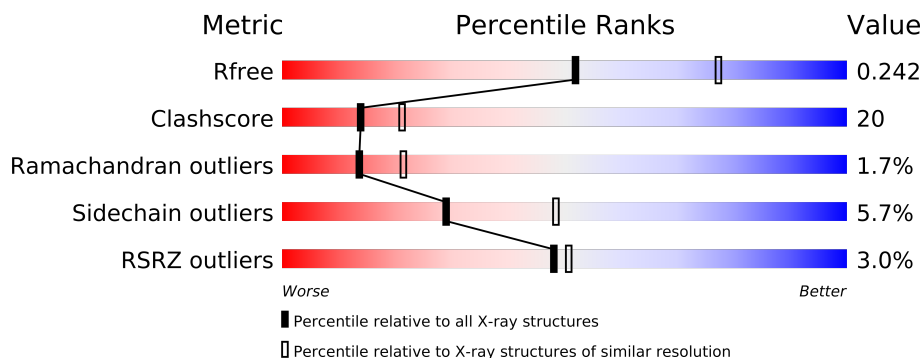
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.54 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))
$R_{free}$	66092	3413 (2.58-2.50)
Clashscore	79885	4284 (2.58-2.50)
Ramachandran outliers	78287	4193 (2.58-2.50)
Sidechain outliers	78261	4195 (2.58-2.50)
RSRZ outliers	66119	3414 (2.58-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	148	
1	B	148	
1	C	148	
1	D	148	
1	E	148	
1	F	148	
1	G	148	
1	H	148	
1	I	148	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11026 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Putative tautomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	145	Total	C	N	O	S	0	0	0
			1179	746	204	226	3			
1	B	145	Total	C	N	O	S	0	0	0
			1190	753	204	230	3			
1	C	146	Total	C	N	O	S	0	0	0
			1196	756	205	232	3			
1	D	145	Total	C	N	O	S	0	0	0
			1182	749	204	226	3			
1	E	145	Total	C	N	O	S	0	0	0
			1154	732	199	220	3			
1	F	145	Total	C	N	O	S	0	0	0
			1186	751	204	228	3			
1	G	145	Total	C	N	O	S	0	0	0
			1171	740	203	225	3			
1	H	140	Total	C	N	O	S	0	0	0
			1122	712	197	210	3			
1	I	145	Total	C	N	O	S	0	0	0
			1166	740	203	220	3			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	53	Total	O	0	0
			53	53		
2	B	48	Total	O	0	0
			48	48		
2	C	50	Total	O	0	0
			50	50		
2	D	43	Total	O	0	0
			43	43		
2	E	55	Total	O	0	0
			55	55		

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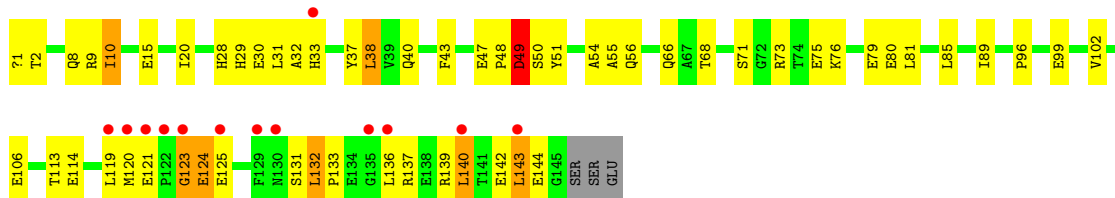
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	F	59	Total 59	O 59	0	0
2	G	49	Total 49	O 49	0	0
2	H	49	Total 49	O 49	0	0
2	I	74	Total 74	O 74	0	0

### 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.

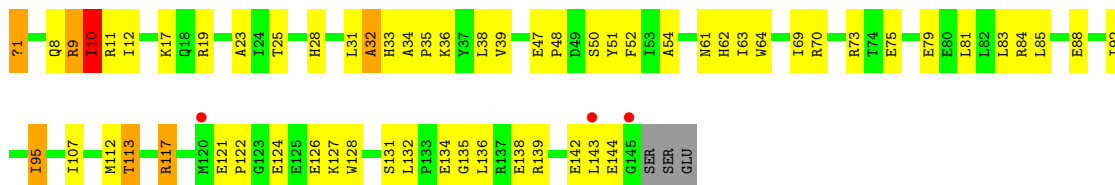
#### • Molecule 1: Putative tautomerase

Chain A: 



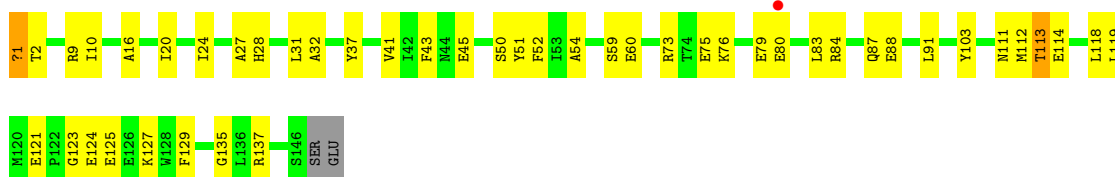
#### • Molecule 1: Putative tautomerase

Chain B: 



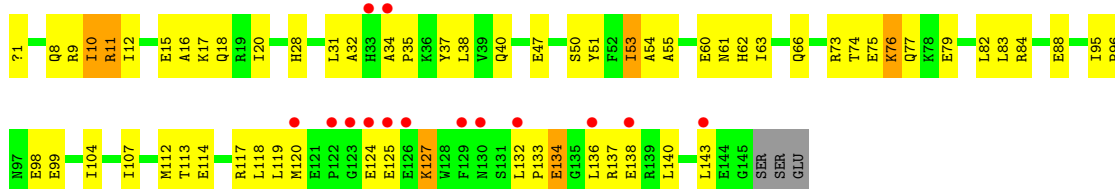
#### • Molecule 1: Putative tautomerase

Chain C: 



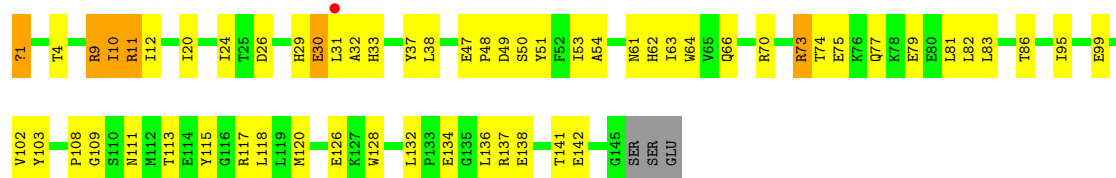
#### • Molecule 1: Putative tautomerase

Chain D: 



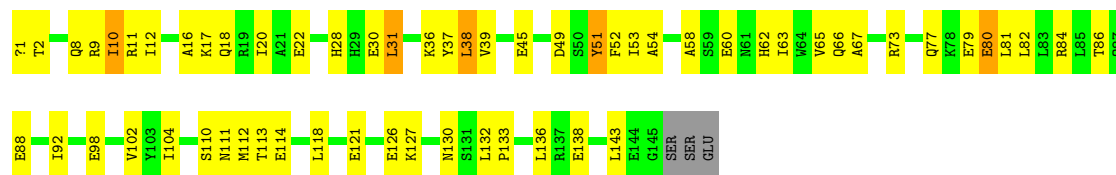
- Molecule 1: Putative tautomerase

Chain E:



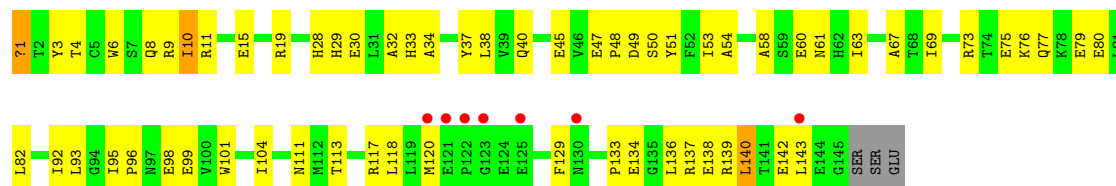
- Molecule 1: Putative tautomerase

Chain F:



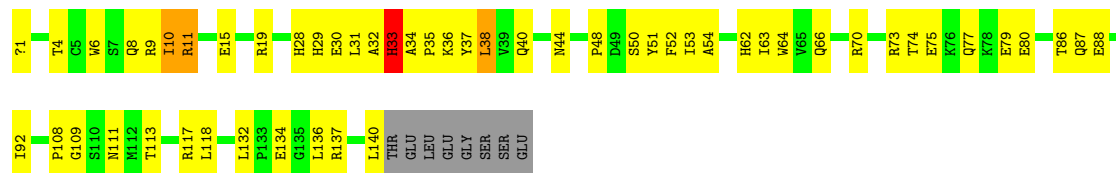
- Molecule 1: Putative tautomerase

Chain G:



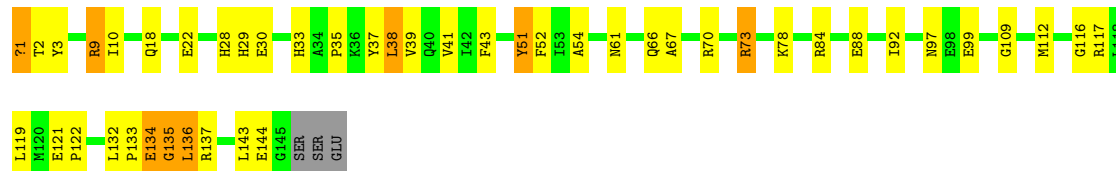
- Molecule 1: Putative tautomerase

Chain H:



- Molecule 1: Putative tautomerase

Chain I:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.10Å 131.00Å 179.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.00 – 2.54 45.21 – 2.02	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.00-2.54) 99.9 (45.21-2.02)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.32 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.223 , 0.243 0.224 , 0.242	Depositor DCC
$R_{free}$ test set	2315 reflections (5.38%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.8	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 32.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	8 of 89205 reflections (0.009%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	11026	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	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 63.44 % 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 9.6669e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PR4

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	0.61	0/1191	0.67	0/1616
1	B	0.57	0/1202	0.61	0/1630
1	C	0.41	0/1208	0.59	0/1638
1	D	0.48	0/1194	0.62	0/1620
1	E	0.58	1/1166 (0.1%)	0.69	1/1587 (0.1%)
1	F	0.49	0/1198	0.63	0/1625
1	G	0.48	0/1183	0.65	0/1607
1	H	0.44	0/1134	0.64	0/1543
1	I	0.64	2/1178 (0.2%)	0.68	0/1601
All	All	0.53	3/10654 (0.0%)	0.64	1/14467 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
1	E	0	1
1	G	0	1
1	I	0	1
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	51	TYR	CD2-CE2	-5.89	1.30	1.39
1	I	51	TYR	CD1-CE1	-5.04	1.31	1.39
1	E	115	TYR	CD1-CE1	-5.02	1.31	1.39



All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	73	ARG	NE-CZ-NH2	-5.62	117.49	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	1	PR4	Mainchain
1	C	1	PR4	Mainchain
1	E	1	PR4	Mainchain
1	G	1	PR4	Mainchain
1	I	1	PR4	Mainchain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1179	0	1150	46	0
1	B	1190	0	1167	54	0
1	C	1196	0	1172	47	0
1	D	1182	0	1159	54	0
1	E	1154	0	1107	52	0
1	F	1186	0	1163	49	0
1	G	1171	0	1127	52	0
1	H	1122	0	1083	47	0
1	I	1166	0	1136	47	0
2	A	53	0	0	10	0
2	B	48	0	0	10	0
2	C	50	0	0	11	0
2	D	43	0	0	11	0
2	E	55	0	0	7	0
2	F	59	0	0	8	0
2	G	49	0	0	10	0
2	H	49	0	0	9	0
2	I	74	0	0	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11026	0	10264	413	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 20.

All (413) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:4:THR:HB	2:H:504:HOH:O	1.32	1.29
1:F:1:PR4:HABA	1:F:112:MET:HE3	1.32	1.10
1:A:113:THR:HG22	2:C:161:HOH:O	1.58	1.03
1:C:1:PR4:HDA	1:C:1:PR4:CAD	1.81	1.03
1:G:9:ARG:HG3	1:G:10:ILE:HG12	1.43	1.00
1:E:11:ARG:HD2	1:E:61:ASN:OD1	1.63	0.99
1:C:1:PR4:HG	1:C:28:HIS:NE2	1.78	0.98
1:E:126:GLU:HB2	2:E:462:HOH:O	1.64	0.96
1:C:125:GLU:HG2	2:C:305:HOH:O	1.67	0.93
1:E:1:PR4:HAC	1:E:70:ARG:HD2	1.51	0.93
1:H:19:ARG:HD2	2:H:156:HOH:O	1.70	0.92
1:C:1:PR4:HDA	1:C:1:PR4:OAG	1.68	0.91
1:I:1:PR4:HABA	1:I:112:MET:HE3	1.53	0.91
1:I:99:GLU:HG3	2:I:445:HOH:O	1.72	0.90
1:D:34:ALA:HB1	2:D:310:HOH:O	1.73	0.88
1:C:10:ILE:HD11	2:C:390:HOH:O	1.73	0.88
1:I:66:GLN:HG2	2:I:367:HOH:O	1.73	0.86
1:E:49:ASP:CB	2:E:152:HOH:O	2.22	0.85
1:B:117:ARG:HD3	2:B:150:HOH:O	1.75	0.85
1:A:68:THR:HB	2:A:553:HOH:O	1.74	0.84
1:G:117:ARG:HD2	2:G:303:HOH:O	1.77	0.84
1:C:1:PR4:OAE	2:C:166:HOH:O	1.93	0.83
1:B:11:ARG:HH21	1:B:95:ILE:HD13	1.43	0.83
1:A:43:PHE:HB2	1:C:50:SER:HB3	1.61	0.82
1:A:2:THR:HB	2:A:553:HOH:O	1.80	0.81
1:A:8:GLN:HG3	2:A:393:HOH:O	1.80	0.81
1:C:1:PR4:CD	1:C:1:PR4:OAG	2.30	0.80
1:F:84:ARG:HD3	2:F:289:HOH:O	1.80	0.80
1:C:1:PR4:CD	1:C:1:PR4:CAD	2.60	0.79
1:F:82:LEU:HD21	1:F:104:ILE:HG12	1.62	0.79
1:A:48:PRO:O	1:A:49:ASP:CB	2.30	0.79
1:F:63:ILE:HD11	2:F:401:HOH:O	1.81	0.79
1:A:102:VAL:H	1:B:113:THR:HG22	1.45	0.79
1:A:66:GLN:HG2	2:A:298:HOH:O	1.84	0.78
1:D:66:GLN:HG3	2:D:321:HOH:O	1.83	0.78

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:84:ARG:HD3	2:D:347:HOH:O	1.84	0.78
1:C:73:ARG:HG2	2:C:291:HOH:O	1.83	0.78
1:B:126:GLU:HB3	2:B:159:HOH:O	1.84	0.77
1:A:48:PRO:O	1:A:49:ASP:HB3	1.84	0.77
1:C:1:PR4:HDA	1:C:1:PR4:OAE	1.85	0.77
1:C:1:PR4:HAC	1:C:112:MET:HE1	1.67	0.76
1:A:120:MET:HB3	1:A:124:GLU:OE1	1.85	0.76
1:B:1:PR4:HDA	1:B:1:PR4:CAD	2.15	0.76
1:H:66:GLN:HG3	2:H:452:HOH:O	1.85	0.76
1:G:34:ALA:HB2	2:G:151:HOH:O	1.85	0.76
1:D:34:ALA:CB	2:D:310:HOH:O	2.32	0.75
1:C:73:ARG:CG	2:C:291:HOH:O	2.34	0.75
1:G:32:ALA:HB3	1:G:73:ARG:HH12	1.50	0.74
1:E:30:GLU:HG3	2:E:419:HOH:O	1.87	0.74
1:D:113:THR:HG22	2:D:315:HOH:O	1.88	0.74
1:I:1:PR4:HABA	1:I:112:MET:CE	2.17	0.74
1:E:74:THR:H	1:E:77:GLN:HE21	1.33	0.74
1:B:61:ASN:HB3	1:B:95:ILE:HG12	1.69	0.74
1:D:117:ARG:HH21	1:D:132:LEU:HA	1.52	0.73
1:E:74:THR:H	1:E:77:GLN:NE2	1.86	0.72
1:H:11:ARG:HD3	1:H:11:ARG:O	1.89	0.72
1:H:140:LEU:CB	2:H:149:HOH:O	2.37	0.72
1:A:15:GLU:HB2	2:A:541:HOH:O	1.89	0.72
1:H:113:THR:HG22	1:H:118:LEU:HD23	1.70	0.72
1:H:117:ARG:HG3	1:H:132:LEU:HD23	1.71	0.72
1:B:113:THR:HG23	2:B:153:HOH:O	1.87	0.72
1:B:47:GLU:HG3	2:B:521:HOH:O	1.89	0.72
1:D:74:THR:HG23	1:D:77:GLN:H	1.55	0.71
1:F:1:PR4:CAB	1:F:112:MET:HE3	2.18	0.71
1:H:32:ALA:O	1:H:34:ALA:N	2.24	0.70
1:E:138:GLU:HA	1:E:141:THR:HG22	1.74	0.69
1:B:124:GLU:HB3	1:B:127:LYS:HD2	1.72	0.69
1:G:32:ALA:HB3	1:G:73:ARG:NH1	2.08	0.69
1:D:88:GLU:HG2	2:D:479:HOH:O	1.93	0.69
1:F:1:PR4:HABA	1:F:112:MET:CE	2.17	0.69
1:G:54:ALA:HB2	1:H:37:TYR:HA	1.74	0.69
1:B:117:ARG:NH1	1:B:117:ARG:HB2	2.08	0.68
1:D:37:TYR:HA	1:F:54:ALA:HB2	1.76	0.68
1:C:27:ALA:O	1:C:31:LEU:HG	1.92	0.68
1:D:9:ARG:O	1:D:10:ILE:HB	1.92	0.68
1:D:134:GLU:HG2	1:D:137:ARG:HH21	1.58	0.68
2:D:403:HOH:O	1:H:48:PRO:HD3	1.93	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:37:TYR:HB3	1:I:143:LEU:HD12	1.76	0.67
1:B:31:LEU:HD11	1:B:84:ARG:HD3	1.75	0.67
1:B:9:ARG:O	1:B:10:ILE:HG12	1.95	0.67
1:G:63:ILE:HD11	1:G:93:LEU:HD12	1.78	0.66
1:H:9:ARG:O	1:H:10:ILE:HG12	1.96	0.66
1:E:47:GLU:O	1:E:50:SER:HB2	1.96	0.66
1:C:1:PR4:HG	1:C:28:HIS:CE1	2.31	0.65
2:A:553:HOH:O	1:C:103:TYR:HE2	1.80	0.65
1:H:74:THR:H	1:H:77:GLN:NE2	1.95	0.65
1:I:134:GLU:O	1:I:136:LEU:N	2.30	0.65
1:D:96:PRO:HD2	1:D:99:GLU:HG3	1.78	0.65
1:B:64:TRP:CH2	1:C:2:THR:HG21	2.31	0.65
1:B:73:ARG:HB2	2:B:156:HOH:O	1.97	0.65
1:E:1:PR4:CAD	1:E:70:ARG:HD3	2.27	0.64
1:F:66:GLN:HG2	2:F:446:HOH:O	1.97	0.64
1:C:1:PR4:HDA	1:C:1:PR4:HOAG	1.63	0.64
1:B:1:PR4:HAC	1:B:70:ARG:HD2	1.81	0.62
1:B:50:SER:HB3	1:C:43:PHE:HB2	1.80	0.62
1:D:32:ALA:HB3	1:D:73:ARG:HH21	1.64	0.62
1:D:47:GLU:O	1:D:50:SER:HB2	2.00	0.62
1:H:75:GLU:O	1:H:79:GLU:HG3	1.99	0.62
1:H:80:GLU:HB3	2:H:506:HOH:O	1.99	0.62
1:I:122:PRO:HD3	2:I:379:HOH:O	1.99	0.62
1:H:11:ARG:HD3	1:H:11:ARG:C	2.19	0.61
1:D:125:GLU:CB	2:D:292:HOH:O	2.48	0.61
1:H:8:GLN:HE21	1:H:48:PRO:HA	1.66	0.61
1:D:120:MET:HG3	1:D:124:GLU:HB2	1.81	0.61
1:D:127:LYS:HZ3	1:D:127:LYS:HB2	1.66	0.61
1:D:76:LYS:HE2	1:D:76:LYS:HA	1.81	0.61
1:D:117:ARG:NH2	1:D:133:PRO:HD3	2.16	0.60
1:G:113:THR:HG22	1:G:118:LEU:HD23	1.83	0.60
1:G:15:GLU:HG2	2:G:149:HOH:O	2.01	0.60
1:F:51:TYR:OH	1:F:62:HIS:HD2	1.84	0.60
1:D:82:LEU:HD21	1:D:104:ILE:HG13	1.81	0.60
1:F:88:GLU:O	1:F:92:ILE:HG13	2.01	0.60
1:B:121:GLU:HG3	1:B:122:PRO:HD2	1.83	0.60
1:C:10:ILE:CD1	2:C:390:HOH:O	2.39	0.60
1:A:47:GLU:O	1:A:50:SER:HB2	2.02	0.60
1:D:66:GLN:CG	2:D:321:HOH:O	2.45	0.59
1:A:37:TYR:HA	1:C:54:ALA:HB2	1.84	0.59
1:B:33:HIS:HB3	2:B:498:HOH:O	2.00	0.59
1:G:82:LEU:HD21	1:G:104:ILE:HG13	1.83	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:9:ARG:O	1:G:10:ILE:HB	2.03	0.59
1:A:96:PRO:HG2	1:A:99:GLU:HG3	1.85	0.59
1:D:54:ALA:O	1:D:55:ALA:HB3	2.02	0.59
1:C:9:ARG:O	1:C:10:ILE:HD12	2.02	0.59
1:C:124:GLU:HG3	1:C:127:LYS:HB3	1.85	0.59
1:E:33:HIS:CE1	2:E:419:HOH:O	2.55	0.59
1:E:74:THR:HG23	1:E:77:GLN:HE21	1.67	0.59
1:A:28:HIS:O	1:A:31:LEU:HB3	2.02	0.59
1:E:1:PR4:CAD	1:E:70:ARG:CD	2.80	0.58
1:B:61:ASN:CG	1:B:95:ILE:HD11	2.23	0.58
1:I:73:ARG:HG2	2:I:153:HOH:O	2.03	0.58
1:B:11:ARG:NH2	1:B:95:ILE:HD13	2.15	0.58
1:E:30:GLU:CG	2:E:419:HOH:O	2.48	0.58
1:H:134:GLU:HA	1:H:137:ARG:HE	1.69	0.58
1:D:1:PR4:HG	1:D:28:HIS:CE1	2.39	0.57
1:E:64:TRP:CH2	1:F:2:THR:HG21	2.39	0.57
1:A:143:LEU:HD13	1:A:144:GLU:N	2.18	0.57
1:B:23:ALA:HB1	1:B:88:GLU:OE2	2.05	0.57
1:E:1:PR4:OAE	1:E:70:ARG:HD3	2.05	0.57
1:D:117:ARG:HD2	2:D:432:HOH:O	2.05	0.57
1:G:11:ARG:HG3	1:G:60:GLU:O	2.05	0.57
1:E:103:TYR:CD1	1:F:112:MET:HG2	2.39	0.57
1:E:1:PR4:CAD	1:E:1:PR4:HDA	2.34	0.57
1:B:117:ARG:CD	2:B:150:HOH:O	2.45	0.57
1:G:34:ALA:CB	2:G:151:HOH:O	2.48	0.56
1:I:121:GLU:CB	2:I:483:HOH:O	2.52	0.56
1:F:80:GLU:HG3	1:F:81:LEU:N	2.20	0.56
1:B:19:ARG:HB3	1:B:92:ILE:HD13	1.88	0.56
1:G:47:GLU:O	1:G:50:SER:HB2	2.05	0.56
1:I:97:ASN:N	1:I:97:ASN:HD22	2.02	0.56
1:G:96:PRO:HG2	1:G:99:GLU:HG2	1.87	0.56
1:E:1:PR4:CAC	1:E:70:ARG:HD2	2.31	0.56
1:E:75:GLU:O	1:E:79:GLU:HG3	2.05	0.56
1:A:29:HIS:C	1:A:31:LEU:H	2.08	0.55
1:I:134:GLU:O	1:I:137:ARG:N	2.36	0.55
1:G:8:GLN:HG3	2:G:302:HOH:O	2.05	0.55
1:E:31:LEU:HD11	1:E:81:LEU:HD22	1.87	0.55
1:G:49:ASP:HA	2:G:407:HOH:O	2.06	0.55
1:I:99:GLU:CG	2:I:445:HOH:O	2.40	0.55
1:D:107:ILE:HD11	1:D:112:MET:HE2	1.89	0.55
1:H:54:ALA:HB2	1:I:37:TYR:HA	1.89	0.55
1:H:86:THR:HG21	1:I:116:GLY:HA2	1.89	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:134:GLU:CB	1:H:137:ARG:HH21	2.20	0.54
1:E:95:ILE:HD12	1:E:99:GLU:HB2	1.89	0.54
1:B:117:ARG:HB2	1:B:117:ARG:HH11	1.72	0.54
1:D:54:ALA:HB2	1:E:37:TYR:HA	1.89	0.54
1:D:75:GLU:HG2	1:D:79:GLU:OE2	2.06	0.54
1:I:35:PRO:HD2	1:I:38:LEU:HD22	1.90	0.54
1:G:63:ILE:CD1	1:G:93:LEU:HD12	2.37	0.54
1:B:12:ILE:HG22	1:B:17:LYS:HG3	1.88	0.54
1:H:35:PRO:HD2	1:H:38:LEU:HD22	1.90	0.54
1:H:62:HIS:O	1:H:63:ILE:HD13	2.07	0.54
1:G:1:PR4:HG	1:G:28:HIS:CE1	2.43	0.54
1:H:1:PR4:HG	1:H:28:HIS:CD2	2.42	0.54
1:B:52:PHE:HB2	1:C:41:VAL:HB	1.90	0.54
1:F:18:GLN:O	1:F:22:GLU:HG3	2.08	0.54
1:C:73:ARG:HB2	2:C:291:HOH:O	2.08	0.54
1:G:1:PR4:CAD	1:G:1:PR4:HDA	2.37	0.54
1:A:9:ARG:O	1:A:10:ILE:HB	2.08	0.54
1:C:73:ARG:CB	2:C:291:HOH:O	2.55	0.53
1:B:1:PR4:HG	1:B:28:HIS:CE1	2.44	0.53
1:B:47:GLU:HG3	1:B:48:PRO:HD2	1.90	0.53
1:B:84:ARG:NH1	2:B:517:HOH:O	2.40	0.53
1:I:132:LEU:C	1:I:133:PRO:O	2.40	0.53
1:C:1:PR4:CG	1:C:28:HIS:NE2	2.64	0.53
1:D:114:GLU:HB2	1:D:119:LEU:HD21	1.91	0.53
1:G:77:GLN:HA	2:G:499:HOH:O	2.06	0.53
1:B:132:LEU:HD22	1:B:136:LEU:HD23	1.90	0.53
1:F:9:ARG:O	1:F:10:ILE:HB	2.07	0.53
1:D:1:PR4:HG	1:D:28:HIS:NE2	2.24	0.53
1:E:12:ILE:CB	2:E:151:HOH:O	2.57	0.53
1:I:30:GLU:HG3	1:I:84:ARG:HH12	1.74	0.53
1:G:75:GLU:O	1:G:79:GLU:HG3	2.09	0.53
1:A:71:SER:HB3	1:A:106:GLU:HB3	1.91	0.53
1:B:1:PR4:CAC	1:B:70:ARG:HD2	2.39	0.52
1:H:30:GLU:CB	2:H:490:HOH:O	2.57	0.52
1:B:1:PR4:CD	1:B:1:PR4:CAD	2.87	0.52
1:C:84:ARG:O	1:C:88:GLU:HG3	2.10	0.52
1:I:1:PR4:CG	1:I:39:VAL:HG22	2.40	0.52
1:B:135:GLY:O	1:B:139:ARG:HD3	2.09	0.52
1:H:29:HIS:HB2	2:H:557:HOH:O	2.10	0.52
1:D:134:GLU:HG2	1:D:137:ARG:NH2	2.24	0.52
1:I:30:GLU:HG3	1:I:84:ARG:NH1	2.25	0.52
1:G:61:ASN:HB3	1:G:95:ILE:HG21	1.91	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:48:PRO:O	1:G:49:ASP:HB2	2.10	0.51
1:D:8:GLN:O	1:D:11:ARG:HG2	2.09	0.51
1:F:133:PRO:HD2	1:F:136:LEU:HD23	1.92	0.51
1:F:12:ILE:HG22	1:F:17:LYS:HG3	1.91	0.51
1:C:9:ARG:HA	1:C:45:GLU:HB3	1.93	0.51
1:B:31:LEU:O	1:B:32:ALA:HB2	2.11	0.51
1:I:88:GLU:O	1:I:92:ILE:HG13	2.10	0.51
1:F:132:LEU:HD23	1:F:136:LEU:HG	1.92	0.51
1:D:134:GLU:O	1:D:138:GLU:HG3	2.10	0.51
1:I:133:PRO:O	1:I:134:GLU:O	2.29	0.51
1:E:54:ALA:HB2	1:F:37:TYR:HA	1.92	0.51
1:I:132:LEU:HB3	1:I:133:PRO:HD2	1.92	0.51
1:D:28:HIS:O	1:D:31:LEU:HB2	2.11	0.51
1:F:77:GLN:O	1:F:80:GLU:HG3	2.11	0.51
1:F:127:LYS:HA	2:F:546:HOH:O	2.10	0.51
1:I:84:ARG:O	1:I:88:GLU:HG3	2.10	0.50
1:H:80:GLU:CB	2:H:506:HOH:O	2.59	0.50
1:G:37:TYR:CE1	1:G:38:LEU:HD13	2.46	0.50
1:C:31:LEU:HD23	1:C:31:LEU:N	2.26	0.50
1:G:15:GLU:HB3	2:G:383:HOH:O	2.10	0.50
1:F:9:ARG:HA	1:F:45:GLU:HB3	1.94	0.50
1:E:48:PRO:O	1:E:49:ASP:CB	2.58	0.50
1:A:1:PR4:HG	1:A:28:HIS:CE1	2.47	0.50
1:C:113:THR:HG22	1:C:113:THR:O	2.11	0.50
1:A:48:PRO:O	1:A:49:ASP:CG	2.50	0.50
1:G:1:PR4:OAF	1:G:69:ILE:HA	2.12	0.50
1:F:65:VAL:HB	1:F:102:VAL:HG22	1.94	0.50
1:A:48:PRO:O	1:A:49:ASP:OD1	2.29	0.49
1:B:1:PR4:HDA	1:B:1:PR4:OAE	2.12	0.49
1:G:136:LEU:HA	1:G:139:ARG:HH21	1.77	0.49
1:D:9:ARG:HG3	1:D:10:ILE:HG12	1.94	0.49
1:I:35:PRO:O	1:I:38:LEU:HB2	2.12	0.49
1:G:98:GLU:OE2	1:H:136:LEU:HD13	2.12	0.49
1:E:62:HIS:O	1:E:63:ILE:HD13	2.11	0.49
1:D:98:GLU:OE2	1:E:136:LEU:HD13	2.13	0.49
1:I:117:ARG:HH11	1:I:117:ARG:HG3	1.77	0.49
1:D:61:ASN:HB3	1:D:95:ILE:HG21	1.94	0.49
1:C:119:LEU:N	1:C:119:LEU:HD22	2.28	0.49
1:A:136:LEU:HA	1:A:139:ARG:HE	1.78	0.49
1:D:1:PR4:HG	1:D:28:HIS:CD2	2.48	0.49
1:D:11:ARG:HG3	1:D:60:GLU:O	2.13	0.49
1:E:4:THR:OG1	1:E:66:GLN:HB3	2.12	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:108:PRO:HG2	1:H:111:ASN:CG	2.32	0.49
1:G:129:PHE:CE2	1:G:137:ARG:HD2	2.47	0.49
1:D:118:LEU:HD13	1:F:79:GLU:HG2	1.95	0.48
1:H:50:SER:HB3	1:I:43:PHE:HB2	1.95	0.48
1:F:30:GLU:HG3	1:F:84:ARG:NH1	2.28	0.48
1:D:74:THR:HG22	1:D:77:GLN:OE1	2.13	0.48
1:I:134:GLU:O	1:I:135:GLY:C	2.51	0.48
1:F:110:SER:C	1:F:111:ASN:HD22	2.16	0.48
1:A:68:THR:O	2:A:553:HOH:O	2.20	0.48
1:H:74:THR:HG23	1:H:77:GLN:NE2	2.27	0.48
1:B:33:HIS:HA	2:B:152:HOH:O	2.12	0.48
1:E:32:ALA:O	1:E:33:HIS:HB2	2.13	0.48
1:F:1:PR4:CG	1:F:39:VAL:HG22	2.43	0.48
1:H:64:TRP:CH2	1:I:2:THR:HG21	2.49	0.48
1:I:18:GLN:HG3	2:I:561:HOH:O	2.12	0.48
1:H:52:PHE:HB2	1:I:41:VAL:HB	1.96	0.48
1:B:134:GLU:O	1:B:138:GLU:HG3	2.14	0.48
1:G:19:ARG:CZ	1:G:92:ILE:HG23	2.44	0.48
1:G:34:ALA:CB	2:G:389:HOH:O	2.62	0.47
1:C:20:ILE:O	1:C:24:ILE:HG13	2.14	0.47
1:B:25:THR:HG23	1:B:39:VAL:HB	1.95	0.47
1:E:134:GLU:HA	1:E:137:ARG:CZ	2.44	0.47
1:C:79:GLU:O	1:C:83:LEU:HG	2.15	0.47
1:B:128:TRP:O	1:B:131:SER:HB3	2.14	0.47
1:C:16:ALA:O	1:C:20:ILE:HG13	2.15	0.47
1:F:1:PR4:HAB	1:F:114:GLU:OE2	2.15	0.47
1:E:108:PRO:O	1:E:111:ASN:HB2	2.15	0.47
1:C:87:GLN:O	1:C:91:LEU:HD23	2.15	0.47
1:B:75:GLU:HG3	1:B:79:GLU:OE2	2.14	0.47
1:I:70:ARG:HB2	1:I:73:ARG:NH2	2.30	0.47
1:G:134:GLU:HA	1:G:137:ARG:CZ	2.44	0.47
1:E:30:GLU:CD	2:E:419:HOH:O	2.52	0.47
1:D:62:HIS:O	1:D:63:ILE:HD13	2.14	0.46
1:E:138:GLU:HA	1:E:141:THR:CG2	2.42	0.46
1:E:53:ILE:O	1:E:54:ALA:HB3	2.16	0.46
1:A:81:LEU:O	1:A:85:LEU:HG	2.15	0.46
1:E:117:ARG:HG3	1:E:132:LEU:HD23	1.98	0.46
1:E:109:GLY:HA3	1:E:120:MET:O	2.15	0.46
1:G:134:GLU:O	1:G:138:GLU:HG2	2.15	0.46
1:E:134:GLU:HA	1:E:137:ARG:NH1	2.31	0.46
1:E:74:THR:HG23	1:E:77:GLN:NE2	2.31	0.46
1:C:114:GLU:HG3	1:C:119:LEU:HD21	1.96	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:32:ALA:C	1:H:34:ALA:N	2.70	0.46
1:B:31:LEU:HD13	1:B:81:LEU:HA	1.98	0.46
1:A:132:LEU:O	1:A:133:PRO:C	2.53	0.46
1:A:76:LYS:O	1:A:80:GLU:HG3	2.16	0.46
1:C:76:LYS:O	1:C:80:GLU:HG3	2.15	0.46
1:H:113:THR:CG2	1:H:118:LEU:HD23	2.44	0.46
1:F:77:GLN:HA	1:F:80:GLU:HG2	1.98	0.46
1:B:54:ALA:HB2	1:C:37:TYR:HA	1.98	0.46
1:H:53:ILE:O	1:H:54:ALA:HB3	2.17	0.45
1:A:37:TYR:CE1	1:A:38:LEU:HD13	2.52	0.45
1:G:29:HIS:O	1:G:33:HIS:HA	2.16	0.45
1:A:137:ARG:HH11	1:A:137:ARG:HG2	1.81	0.45
1:F:16:ALA:O	1:F:20:ILE:HG13	2.15	0.45
1:D:95:ILE:O	1:I:9:ARG:NH2	2.49	0.45
1:D:83:LEU:HD21	1:E:118:LEU:CD2	2.46	0.45
1:B:61:ASN:HB3	1:B:95:ILE:CG1	2.43	0.45
1:I:132:LEU:O	1:I:133:PRO:C	2.53	0.45
1:A:8:GLN:NE2	2:A:533:HOH:O	2.50	0.45
1:A:75:GLU:O	1:A:79:GLU:HG3	2.16	0.45
1:E:11:ARG:HH11	1:E:61:ASN:HB3	1.81	0.45
1:F:126:GLU:HB3	2:F:402:HOH:O	2.17	0.45
1:D:35:PRO:HG2	1:D:38:LEU:HD22	1.99	0.45
1:G:79:GLU:HB3	1:H:118:LEU:HD11	1.99	0.45
1:B:36:LYS:HB2	1:B:144:GLU:OE2	2.17	0.45
1:I:132:LEU:O	1:I:133:PRO:O	2.35	0.45
1:D:73:ARG:HB2	2:D:341:HOH:O	2.16	0.45
1:I:18:GLN:O	1:I:22:GLU:HG3	2.17	0.45
1:D:12:ILE:HG22	1:D:17:LYS:HG3	1.99	0.45
1:A:8:GLN:CG	2:A:393:HOH:O	2.53	0.45
1:B:1:PR4:OAF	1:B:69:ILE:HA	2.17	0.45
1:B:8:GLN:HG3	2:B:411:HOH:O	2.18	0.45
1:I:29:HIS:O	1:I:33:HIS:HA	2.17	0.45
1:H:32:ALA:O	1:H:33:HIS:C	2.56	0.44
1:I:61:ASN:HB2	2:I:312:HOH:O	2.17	0.44
1:B:62:HIS:O	1:B:63:ILE:HD13	2.17	0.44
1:F:1:PR4:CAB	1:F:112:MET:CE	2.89	0.44
1:E:118:LEU:O	1:E:128:TRP:HD1	2.00	0.44
1:F:53:ILE:HG12	1:F:58:ALA:HB2	2.00	0.44
1:D:16:ALA:O	1:D:20:ILE:HG13	2.18	0.44
1:B:61:ASN:OD1	1:B:95:ILE:HD11	2.17	0.44
1:A:29:HIS:C	1:A:31:LEU:N	2.71	0.44
1:H:108:PRO:O	1:H:111:ASN:HB2	2.18	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:123:GLY:O	1:A:124:GLU:HG3	2.18	0.44
1:E:82:LEU:HD23	1:E:102:VAL:CG1	2.48	0.44
1:G:40:GLN:HA	1:I:52:PHE:O	2.18	0.44
1:D:127:LYS:NZ	1:D:127:LYS:HB2	2.32	0.44
1:I:97:ASN:N	1:I:97:ASN:ND2	2.66	0.44
1:A:33:HIS:CE1	2:A:486:HOH:O	2.71	0.44
1:I:109:GLY:O	1:I:119:LEU:O	2.36	0.44
1:A:32:ALA:HB3	1:A:73:ARG:NH1	2.32	0.44
1:E:138:GLU:O	1:E:142:GLU:HG3	2.17	0.43
1:D:136:LEU:O	1:D:140:LEU:HG	2.18	0.43
1:A:136:LEU:CB	1:A:139:ARG:HH21	2.31	0.43
1:D:113:THR:HG21	1:F:86:THR:OG1	2.18	0.43
1:A:56:GLN:NE2	1:B:143:LEU:HD22	2.33	0.43
1:G:134:GLU:HA	1:G:137:ARG:NH1	2.33	0.43
1:I:1:PR4:HG	1:I:28:HIS:CD2	2.54	0.43
1:G:111:ASN:HD21	1:I:78:LYS:NZ	2.16	0.43
1:H:70:ARG:HG2	1:H:109:GLY:HA2	2.00	0.43
1:G:140:LEU:O	1:G:142:GLU:N	2.47	0.43
1:A:29:HIS:O	1:A:31:LEU:N	2.42	0.43
1:G:133:PRO:HG2	1:G:136:LEU:HB2	2.01	0.42
1:B:34:ALA:HB1	1:B:35:PRO:HD2	2.00	0.42
1:H:74:THR:HG23	1:H:77:GLN:HE21	1.83	0.42
1:G:37:TYR:HA	1:I:54:ALA:HB2	2.02	0.42
1:C:123:GLY:HA2	2:C:305:HOH:O	2.19	0.42
1:G:34:ALA:HB1	2:G:389:HOH:O	2.19	0.42
1:I:18:GLN:CG	2:I:561:HOH:O	2.68	0.42
1:D:40:GLN:HA	1:F:52:PHE:O	2.19	0.42
1:B:11:ARG:HH21	1:B:95:ILE:CD1	2.22	0.42
1:B:70:ARG:HA	1:B:107:ILE:O	2.20	0.42
1:H:28:HIS:O	1:H:31:LEU:HB2	2.19	0.42
1:F:67:ALA:HB3	1:F:104:ILE:HD13	2.01	0.42
1:F:8:GLN:HB3	1:F:11:ARG:HG3	2.02	0.42
1:G:4:THR:HG22	1:G:6:TRP:CD1	2.54	0.42
1:I:3:TYR:CD2	1:I:67:ALA:HB2	2.54	0.42
1:C:1:PR4:CG	1:C:28:HIS:CD2	3.03	0.42
1:A:54:ALA:O	1:A:55:ALA:HB3	2.20	0.42
1:C:1:PR4:HG	1:C:28:HIS:CD2	2.51	0.42
1:E:1:PR4:HAC	1:E:70:ARG:CD	2.37	0.42
1:F:62:HIS:O	1:F:63:ILE:HD13	2.20	0.42
1:A:121:GLU:O	1:A:124:GLU:HB2	2.20	0.42
1:E:79:GLU:HB3	1:F:118:LEU:HD11	2.00	0.42
1:E:138:GLU:CA	1:E:141:THR:HG22	2.46	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:121:GLU:HG3	2:F:435:HOH:O	2.20	0.42
1:B:83:LEU:HD21	1:C:118:LEU:HG	2.01	0.42
1:F:73:ARG:N	1:F:73:ARG:HD2	2.35	0.41
1:G:134:GLU:HG2	1:G:137:ARG:NH2	2.35	0.41
1:E:86:THR:OG1	1:F:113:THR:HG21	2.19	0.41
1:G:101:TRP:CZ3	1:H:40:GLN:HB3	2.55	0.41
1:H:6:TRP:CZ3	1:H:44:ASN:ND2	2.88	0.41
1:E:26:ASP:O	1:E:29:HIS:HB3	2.20	0.41
1:B:70:ARG:HG3	1:B:112:MET:SD	2.61	0.41
1:G:76:LYS:NZ	1:G:80:GLU:OE1	2.52	0.41
1:F:1:PR4:HG	1:F:28:HIS:CD2	2.55	0.41
1:G:9:ARG:HE	1:G:9:ARG:HB2	1.71	0.41
1:D:117:ARG:NH2	1:D:132:LEU:HA	2.25	0.41
1:E:113:THR:HG22	1:E:118:LEU:HD13	2.03	0.41
1:F:36:LYS:C	1:F:38:LEU:H	2.24	0.41
1:F:30:GLU:HG3	1:F:84:ARG:HH12	1.85	0.41
1:E:82:LEU:HD23	1:E:102:VAL:HB	2.03	0.41
1:F:8:GLN:HB3	1:F:11:ARG:CG	2.51	0.41
1:A:20:ILE:HG23	1:A:89:ILE:HG12	2.02	0.41
1:G:53:ILE:HG12	1:G:58:ALA:HB2	2.02	0.41
1:C:129:PHE:CE2	1:C:137:ARG:HD2	2.56	0.41
1:F:51:TYR:OH	1:F:62:HIS:CD2	2.69	0.41
1:H:32:ALA:N	2:H:507:HOH:O	2.54	0.41
1:G:63:ILE:HD11	1:G:93:LEU:CD1	2.49	0.41
1:I:144:GLU:HG3	2:I:472:HOH:O	2.21	0.41
1:A:114:GLU:HG3	1:A:119:LEU:HD11	2.03	0.41
1:D:143:LEU:O	1:D:143:LEU:HD23	2.21	0.41
1:G:96:PRO:HG2	1:G:99:GLU:CG	2.51	0.40
1:C:135:GLY:HA3	2:C:150:HOH:O	2.20	0.40
1:E:20:ILE:O	1:E:24:ILE:HG13	2.20	0.40
1:I:66:GLN:CG	2:I:367:HOH:O	2.49	0.40
1:H:32:ALA:CB	1:H:73:ARG:HH12	2.35	0.40
1:A:96:PRO:HG2	1:A:99:GLU:CG	2.50	0.40
1:H:29:HIS:HD2	1:H:36:LYS:HG3	1.85	0.40
1:A:136:LEU:O	1:A:140:LEU:HD22	2.21	0.40
1:F:60:GLU:HG2	2:F:163:HOH:O	2.21	0.40
1:G:3:TYR:CD2	1:G:67:ALA:HB2	2.56	0.40
1:C:32:ALA:HB1	1:C:73:ARG:CZ	2.52	0.40
1:F:49:ASP:HA	2:F:395:HOH:O	2.21	0.40
1:A:40:GLN:HA	1:C:52:PHE:O	2.21	0.40
1:H:88:GLU:O	1:H:92:ILE:HG13	2.22	0.40
1:D:53:ILE:O	1:D:54:ALA:HB3	2.21	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	143/148 (97%)	132 (92%)	5 (4%)	6 (4%)	4	4
1	B	143/148 (97%)	136 (95%)	4 (3%)	3 (2%)	11	16
1	C	144/148 (97%)	137 (95%)	6 (4%)	1 (1%)	30	49
1	D	143/148 (97%)	137 (96%)	5 (4%)	1 (1%)	30	49
1	E	143/148 (97%)	138 (96%)	3 (2%)	2 (1%)	16	27
1	F	143/148 (97%)	138 (96%)	3 (2%)	2 (1%)	16	27
1	G	143/148 (97%)	136 (95%)	5 (4%)	2 (1%)	16	27
1	H	138/148 (93%)	126 (91%)	10 (7%)	2 (1%)	16	27
1	I	143/148 (97%)	135 (94%)	5 (4%)	3 (2%)	11	16
All	All	1283/1332 (96%)	1215 (95%)	46 (4%)	22 (2%)	14	21

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	GLU
1	B	10	ILE
1	B	32	ALA
1	D	10	ILE
1	E	10	ILE
1	G	10	ILE
1	H	10	ILE
1	H	33	HIS
1	I	134	GLU
1	I	135	GLY
1	A	10	ILE
1	A	49	ASP
1	A	123	GLY
1	B	142	GLU

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Mol	Chain	Res	Type
1	F	10	ILE
1	C	75	GLU
1	F	31	LEU
1	G	140	LEU
1	A	30	GLU
1	E	9	ARG
1	A	131	SER
1	I	10	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	123/129 (95%)	115 (94%)	8 (6%)	24	41
1	B	126/129 (98%)	118 (94%)	8 (6%)	25	42
1	C	127/129 (98%)	121 (95%)	6 (5%)	36	59
1	D	124/129 (96%)	116 (94%)	8 (6%)	24	41
1	E	117/129 (91%)	109 (93%)	8 (7%)	22	38
1	F	125/129 (97%)	117 (94%)	8 (6%)	25	42
1	G	120/129 (93%)	115 (96%)	5 (4%)	40	65
1	H	114/129 (88%)	108 (95%)	6 (5%)	32	53
1	I	120/129 (93%)	115 (96%)	5 (4%)	40	65
All	All	1096/1161 (94%)	1034 (94%)	62 (6%)	29	48

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	LEU
1	A	49	ASP
1	A	51	TYR
1	A	125	GLU
1	A	132	LEU
1	A	140	LEU
1	A	142	GLU

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Mol	Chain	Res	Type
1	A	143	LEU
1	B	9	ARG
1	B	10	ILE
1	B	38	LEU
1	B	51	TYR
1	B	85	LEU
1	B	95	ILE
1	B	113	THR
1	B	117	ARG
1	C	51	TYR
1	C	59	SER
1	C	60	GLU
1	C	111	ASN
1	C	113	THR
1	C	121	GLU
1	D	11	ARG
1	D	15	GLU
1	D	18	GLN
1	D	51	TYR
1	D	53	ILE
1	D	76	LYS
1	D	127	LYS
1	D	134	GLU
1	E	9	ARG
1	E	10	ILE
1	E	11	ARG
1	E	30	GLU
1	E	38	LEU
1	E	51	TYR
1	E	73	ARG
1	E	83	LEU
1	F	31	LEU
1	F	38	LEU
1	F	51	TYR
1	F	80	GLU
1	F	98	GLU
1	F	130	ASN
1	F	138	GLU
1	F	143	LEU
1	G	30	GLU
1	G	45	GLU
1	G	51	TYR

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Mol	Chain	Res	Type
1	G	120	MET
1	G	143	LEU
1	H	11	ARG
1	H	15	GLU
1	H	33	HIS
1	H	38	LEU
1	H	51	TYR
1	H	87	GLN
1	I	9	ARG
1	I	38	LEU
1	I	51	TYR
1	I	73	ARG
1	I	136	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	56	GLN
1	A	62	HIS
1	A	111	ASN
1	B	8	GLN
1	C	18	GLN
1	C	56	GLN
1	C	62	HIS
1	C	130	ASN
1	D	8	GLN
1	D	18	GLN
1	D	66	GLN
1	D	111	ASN
1	D	130	ASN
1	E	77	GLN
1	E	111	ASN
1	F	18	GLN
1	F	56	GLN
1	F	62	HIS
1	F	66	GLN
1	F	111	ASN
1	F	130	ASN
1	G	8	GLN
1	G	18	GLN
1	G	62	HIS

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Mol	Chain	Res	Type
1	G	87	GLN
1	G	111	ASN
1	H	8	GLN
1	H	29	HIS
1	H	77	GLN
1	H	111	ASN
1	I	18	GLN
1	I	62	HIS
1	I	97	ASN
1	I	111	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

9 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
1	PR4	A	1	1	13,13,14	3.73	2 (15%)	14,17,19	2.07	5 (35%)
1	PR4	B	1	1	13,13,14	5.14	3 (23%)	14,17,19	1.10	0
1	PR4	C	1	1	13,13,14	4.53	1 (7%)	14,17,19	1.46	2 (14%)
1	PR4	D	1	1	13,13,14	5.24	2 (15%)	14,17,19	1.63	3 (21%)
1	PR4	E	1	1	13,13,14	5.16	3 (23%)	14,17,19	1.16	0
1	PR4	F	1	1	13,13,14	5.29	2 (15%)	14,17,19	1.09	0
1	PR4	G	1	1	13,13,14	5.32	3 (23%)	14,17,19	1.53	3 (21%)
1	PR4	H	1	1	13,13,14	5.24	1 (7%)	14,17,19	1.48	2 (14%)
1	PR4	I	1	1	13,13,14	4.94	2 (15%)	14,17,19	1.30	2 (14%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral



centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PR4	A	1	1	-	0/8/20/22	0/1/1/1
1	PR4	B	1	1	-	0/8/20/22	0/1/1/1
1	PR4	C	1	1	-	0/8/20/22	0/1/1/1
1	PR4	D	1	1	-	0/8/20/22	0/1/1/1
1	PR4	E	1	1	-	0/8/20/22	0/1/1/1
1	PR4	F	1	1	-	0/8/20/22	0/1/1/1
1	PR4	G	1	1	-	0/8/20/22	0/1/1/1
1	PR4	H	1	1	-	0/8/20/22	0/1/1/1
1	PR4	I	1	1	-	0/8/20/22	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	1	PR4	O-C	18.65	1.24	1.11
1	F	1	PR4	O-C	18.53	1.24	1.11
1	D	1	PR4	O-C	18.53	1.24	1.11
1	G	1	PR4	O-C	18.49	1.24	1.11
1	E	1	PR4	O-C	18.06	1.23	1.11
1	B	1	PR4	O-C	17.95	1.23	1.11
1	I	1	PR4	O-C	17.20	1.23	1.11
1	C	1	PR4	O-C	16.07	1.22	1.11
1	A	1	PR4	O-C	12.55	1.20	1.11
1	G	1	PR4	CA-C	3.89	1.55	1.48
1	F	1	PR4	CA-C	3.16	1.54	1.48
1	E	1	PR4	CAC-CAD	3.07	1.57	1.52
1	A	1	PR4	CB-CA	-2.98	1.49	1.53
1	I	1	PR4	CA-C	2.88	1.53	1.48
1	G	1	PR4	CAC-CAD	2.79	1.56	1.52
1	B	1	PR4	CAC-CAD	2.77	1.56	1.52
1	D	1	PR4	CA-C	2.63	1.53	1.48
1	B	1	PR4	CA-C	2.29	1.52	1.48
1	E	1	PR4	CA-C	2.04	1.52	1.48

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	PR4	OAE-CAD-CAC	5.07	126.83	113.19
1	D	1	PR4	OAF-CAD-CAC	-3.58	113.04	122.12
1	A	1	PR4	OAF-CAD-CAC	-3.02	114.45	122.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	1	PR4	OAF-CAD-CAC	-3.01	114.48	122.12
1	G	1	PR4	OAE-CAD-CAC	2.78	120.67	113.19
1	C	1	PR4	CB-CA-N	2.77	108.19	103.16
1	I	1	PR4	CAC-CAB-N	2.63	119.68	112.75
1	H	1	PR4	CAB-N-CA	-2.61	106.61	113.37
1	D	1	PR4	OAE-CAD-CAC	2.42	119.70	113.19
1	A	1	PR4	OAE-CAD-OAF	-2.38	118.69	124.07
1	H	1	PR4	OAF-CAD-CAC	-2.33	116.21	122.12
1	A	1	PR4	CD-N-CA	2.28	110.97	106.30
1	I	1	PR4	CAB-N-CA	-2.23	107.60	113.37
1	C	1	PR4	CG-CD-N	2.23	108.18	104.13
1	D	1	PR4	CAC-CAB-N	2.22	118.60	112.75
1	A	1	PR4	CAC-CAB-N	2.05	118.17	112.75
1	G	1	PR4	CAC-CAB-N	2.01	118.06	112.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

There are no ligands in this entry.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	145/148 (97%)	0.28	13 (8%) 10 9	13, 31, 77, 79	0
1	B	145/148 (97%)	0.05	3 (2%) 60 64	18, 36, 66, 83	0
1	C	146/148 (98%)	-0.03	1 (0%) 84 87	13, 35, 59, 65	0
1	D	145/148 (97%)	0.34	14 (9%) 8 8	11, 31, 87, 92	0
1	E	145/148 (97%)	-0.18	1 (0%) 84 87	16, 33, 48, 56	0
1	F	145/148 (97%)	-0.07	0 100 100	16, 32, 53, 61	0
1	G	145/148 (97%)	0.15	7 (4%) 29 31	14, 30, 67, 70	0
1	H	140/148 (94%)	-0.17	0 100 100	15, 34, 58, 62	0
1	I	145/148 (97%)	-0.22	0 100 100	10, 28, 56, 62	0
All	All	1301/1332 (97%)	0.02	39 (2%) 48 51	10, 33, 67, 92	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	MET	4.9
1	D	123	GLY	4.8
1	D	136	LEU	4.7
1	G	120	MET	3.9
1	A	140	LEU	3.8
1	G	122	PRO	3.8
1	D	126	GLU	3.8
1	D	129	PHE	3.7
1	G	125	GLU	3.6
1	B	145	GLY	3.5
1	A	119	LEU	3.5
1	D	125	GLU	3.4
1	D	122	PRO	3.4
1	B	143	LEU	3.3
1	A	122	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	136	LEU	3.3
1	A	143	LEU	3.3
1	D	130	ASN	3.2
1	G	143	LEU	2.9
1	E	31	LEU	2.8
1	A	121	GLU	2.8
1	B	120	MET	2.8
1	A	130	ASN	2.7
1	D	143	LEU	2.7
1	A	33	HIS	2.7
1	D	120	MET	2.7
1	D	34	ALA	2.6
1	A	123	GLY	2.4
1	D	124	GLU	2.4
1	G	130	ASN	2.4
1	G	123	GLY	2.3
1	D	33	HIS	2.3
1	C	80	GLU	2.3
1	G	121	GLU	2.3
1	A	135	GLY	2.2
1	D	132	LEU	2.1
1	D	138	GLU	2.1
1	A	125	GLU	2.1
1	A	129	PHE	2.0

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

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	PR4	C	1	13/14	0.27	5.14	28,30,40,41	0
1	PR4	F	1	13/14	0.20	2.22	27,28,34,38	0
1	PR4	E	1	13/14	0.17	1.31	23,25,33,36	0
1	PR4	G	1	13/14	0.20	0.97	29,30,36,36	0
1	PR4	I	1	13/14	0.15	0.88	24,25,28,30	0
1	PR4	A	1	13/14	0.20	0.70	27,29,40,43	0
1	PR4	D	1	13/14	0.17	0.39	30,31,36,36	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	PR4	B	1	13/14	0.15	0.14	27,29,34,38	0
1	PR4	H	1	13/14	0.14	0.00	26,28,33,33	0

### 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.