



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:56 AM GMT

PDB ID : 1FPZ  
Title : CRYSTAL STRUCTURE ANALYSIS OF KINASE ASSOCIATED PHOSPHATASE (KAP) WITH A SUBSTITUTION OF THE CATALYTIC SITE CYSTEINE (CYS140) TO A SERINE  
Authors : Song, H.; Hanlon, N.; Brown, N.R.; Noble, M.E.M.; Johnson, L.N.; Barford, D.  
Deposited on : 2000-09-01  
Resolution : 2.00 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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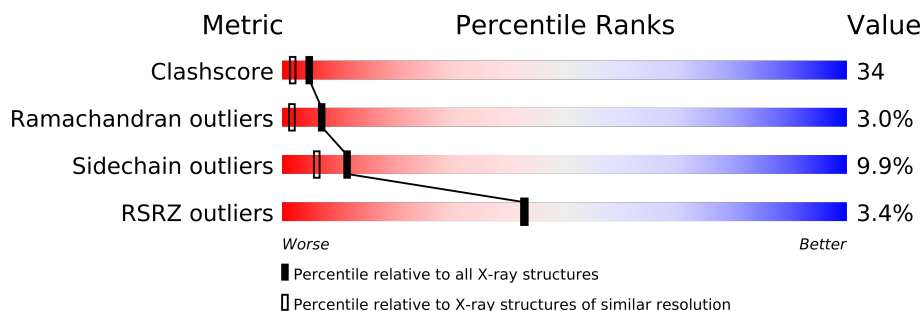
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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	212	
1	B	212	
1	C	212	
1	D	212	
1	E	212	
1	F	212	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9060 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 CYCLIN-DEPENDENT KINASE INHIBITOR 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	176	Total	C	N	O	S	0	0	0
			1381	871	243	255	12			
1	B	176	Total	C	N	O	S	0	0	0
			1381	871	243	255	12			
1	C	178	Total	C	N	O	S	0	0	0
			1399	881	246	260	12			
1	D	178	Total	C	N	O	S	0	0	0
			1399	881	246	260	12			
1	E	176	Total	C	N	O	S	0	0	0
			1381	871	243	255	12			
1	F	178	Total	C	N	O	S	0	0	0
			1399	881	246	260	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	SER	CYS	MUTATION	UNP Q16667
B	140	SER	CYS	MUTATION	UNP Q16667
C	140	SER	CYS	MUTATION	UNP Q16667
D	140	SER	CYS	MUTATION	UNP Q16667
E	140	SER	CYS	MUTATION	UNP Q16667
F	140	SER	CYS	MUTATION	UNP Q16667

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

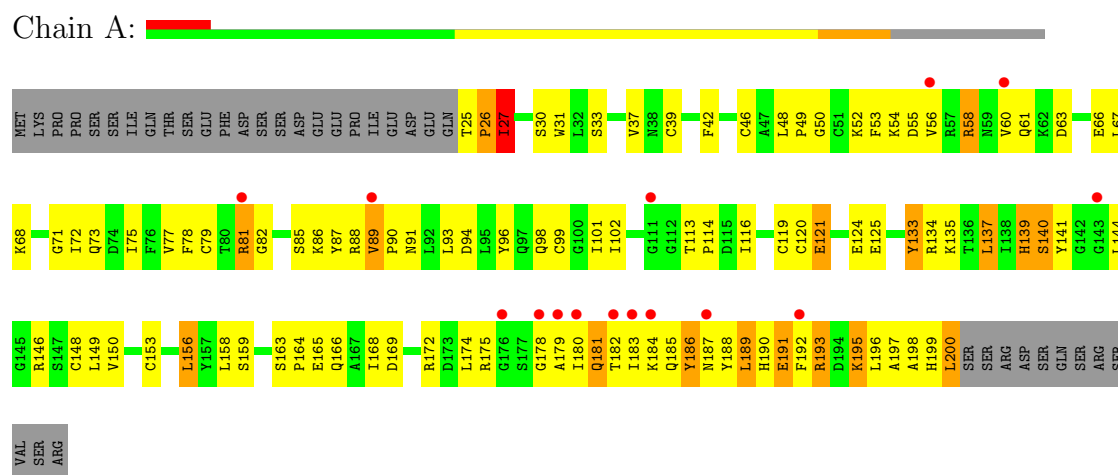
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		
3	B	99	Total	O	0	0
			99	99		
3	C	133	Total	O	0	0
			133	133		
3	D	116	Total	O	0	0
			116	116		
3	E	155	Total	O	0	0
			155	155		
3	F	95	Total	O	0	0
			95	95		

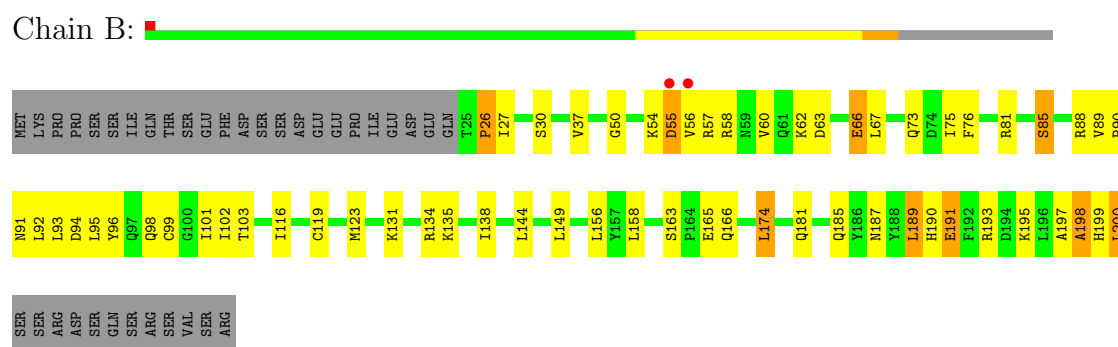
### 3 Residue-property plots

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.

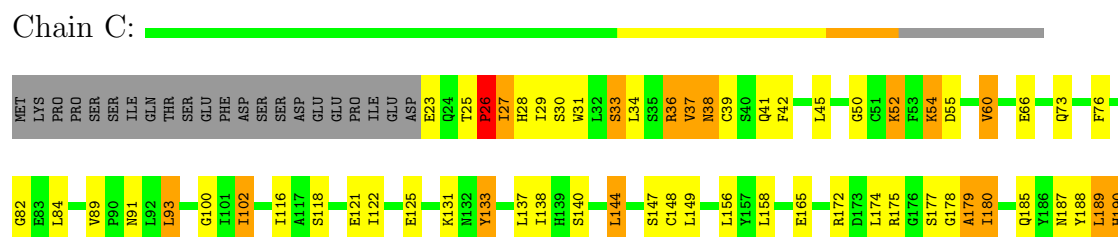
#### • Molecule 1: CYCLIN-DEPENDENT KINASE INHIBITOR 3



#### • Molecule 1: CYCLIN-DEPENDENT KINASE INHIBITOR 3



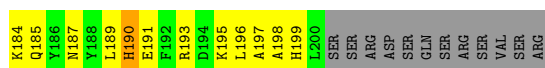
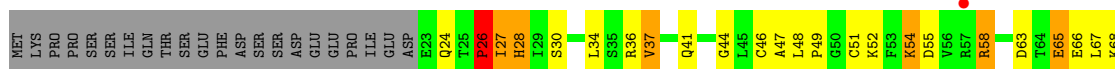
#### • Molecule 1: CYCLIN-DEPENDENT KINASE INHIBITOR 3





• Molecule 1: CYCLIN-DEPENDENT KINASE INHIBITOR 3

Chain D:



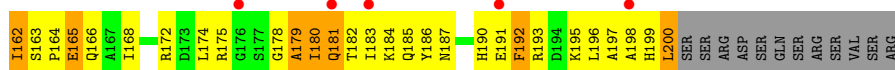
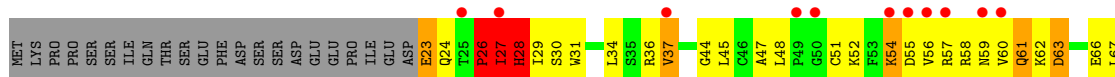
• Molecule 1: CYCLIN-DEPENDENT KINASE INHIBITOR 3

Chain E:



• Molecule 1: CYCLIN-DEPENDENT KINASE INHIBITOR 3

Chain F:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	131.93Å 131.93Å 140.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.00 48.05 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.7 (20.00-2.00) 96.6 (48.05-2.00)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.76 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.202 , 0.253 0.214 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	22.6	Xtriage
Anisotropy	0.416	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 42.7	EDS
Estimated twinning fraction	0.487 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 90052 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9060	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	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 55.54 % 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 3.1165e-05. 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: SO4

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	0.47	0/1407	0.69	0/1902
1	B	0.62	0/1407	0.78	1/1902 (0.1%)
1	C	0.60	0/1425	0.76	2/1926 (0.1%)
1	D	0.57	0/1425	0.77	1/1926 (0.1%)
1	E	0.63	1/1407 (0.1%)	0.79	1/1902 (0.1%)
1	F	0.46	0/1425	0.73	0/1926
All	All	0.56	1/8496 (0.0%)	0.75	5/11484 (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	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	70	CYS	CB-SG	-5.71	1.72	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	144	LEU	CA-CB-CG	5.90	128.88	115.30
1	E	144	LEU	CA-CB-CG	5.52	128.00	115.30
1	D	190	HIS	N-CA-C	-5.47	96.24	111.00
1	C	102	ILE	CG1-CB-CG2	-5.42	99.48	111.40
1	C	190	HIS	N-CA-C	-5.24	96.86	111.00



There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	133	TYR	Sidechain

## 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	1381	0	1389	132	1
1	B	1381	0	1389	68	0
1	C	1399	0	1403	71	0
1	D	1399	0	1403	82	1
1	E	1381	0	1389	79	0
1	F	1399	0	1403	144	0
2	A	5	0	0	1	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	92	0	0	26	0
3	B	99	0	0	5	0
3	C	133	0	0	28	0
3	D	116	0	0	27	0
3	E	155	0	0	26	1
3	F	95	0	0	34	0
All	All	9060	0	8376	575	2

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 34.

All (575) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:27:ILE:HD13	1:B:66:GLU:HG2	1.31	1.11
1:C:29:ILE:HG22	3:C:988:HOH:O	1.56	1.05

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:181:GLN:HE21	1:A:181:GLN:HA	1.27	1.00
1:E:36:ARG:HB2	1:E:36:ARG:HH11	1.28	0.97
1:C:93:LEU:H	1:C:93:LEU:HD22	1.30	0.96
1:D:49:PRO:HD2	3:D:933:HOH:O	1.66	0.95
1:E:124:GLU:HA	3:E:1059:HOH:O	1.65	0.95
1:B:30:SER:HB3	1:B:174:LEU:HD21	1.49	0.93
1:E:152:ALA:O	1:E:156:LEU:HD13	1.69	0.92
1:B:55:ASP:H	1:B:57:ARG:NH2	1.67	0.91
1:D:73:GLN:HE21	1:D:135:LYS:H	1.17	0.91
1:F:85:SER:HA	1:F:90:PRO:HG3	1.53	0.91
1:A:120:CYS:O	1:A:124:GLU:HG3	1.69	0.90
1:A:101:ILE:HD11	3:A:920:HOH:O	1.72	0.90
1:F:163:SER:HB3	1:F:166:GLN:HG3	1.53	0.90
1:E:156:LEU:HG	3:E:948:HOH:O	1.73	0.89
1:B:55:ASP:H	1:B:57:ARG:HH21	0.99	0.88
1:F:179:ALA:O	1:F:181:GLN:HG3	1.74	0.88
1:F:91:ASN:HD22	1:F:95:LEU:HD21	1.37	0.87
1:E:60:VAL:HG23	3:E:909:HOH:O	1.74	0.86
1:F:116:ILE:HG23	1:F:195:LYS:HE2	1.57	0.86
1:E:127:THR:HB	3:E:1059:HOH:O	1.75	0.86
1:A:81:ARG:HG2	3:A:935:HOH:O	1.78	0.83
1:C:93:LEU:HD22	1:C:93:LEU:N	1.93	0.83
1:D:190:HIS:O	1:D:191:GLU:HG2	1.78	0.82
1:F:81:ARG:HB2	1:F:81:ARG:NH1	1.93	0.82
1:C:200:LEU:HD22	1:C:200:LEU:H	1.40	0.82
1:A:93:LEU:HD13	3:A:946:HOH:O	1.80	0.81
1:F:180:ILE:HG12	3:F:926:HOH:O	1.80	0.81
1:F:72:ILE:HD13	1:F:135:LYS:HE2	1.62	0.80
1:B:27:ILE:HG13	1:B:58:ARG:NH2	1.97	0.80
1:F:181:GLN:HB3	3:F:915:HOH:O	1.80	0.80
1:E:36:ARG:HD3	3:E:1033:HOH:O	1.80	0.80
1:F:56:VAL:O	1:F:57:ARG:HD3	1.81	0.79
1:D:198:ALA:HB2	3:D:934:HOH:O	1.83	0.79
1:C:93:LEU:HG	3:C:990:HOH:O	1.81	0.78
1:E:191:GLU:HG3	1:E:195:LYS:HD2	1.62	0.78
1:F:73:GLN:NE2	1:F:135:LYS:H	1.81	0.78
1:F:181:GLN:OE1	1:F:182:THR:HG23	1.84	0.78
1:D:73:GLN:NE2	1:D:135:LYS:H	1.81	0.78
1:D:193:ARG:HD2	3:D:941:HOH:O	1.85	0.77
1:C:190:HIS:O	1:C:191:GLU:HG2	1.85	0.77
1:F:89:VAL:N	1:F:90:PRO:HD3	2.01	0.77
1:A:89:VAL:N	1:A:90:PRO:HD3	2.00	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:68:LYS:HB2	3:E:953:HOH:O	1.85	0.76
1:B:195:LYS:HG2	3:B:966:HOH:O	1.86	0.76
1:C:93:LEU:H	1:C:93:LEU:CD2	1.97	0.75
1:D:198:ALA:HB1	3:D:999:HOH:O	1.85	0.75
1:B:27:ILE:HG13	1:B:58:ARG:HH22	1.53	0.74
1:B:89:VAL:N	1:B:90:PRO:HD3	2.02	0.74
1:F:94:ASP:HB3	3:F:978:HOH:O	1.86	0.74
1:F:91:ASN:ND2	1:F:95:LEU:HD11	2.02	0.74
1:E:168:ILE:HG23	3:E:1043:HOH:O	1.88	0.74
1:C:116:ILE:HG13	3:C:919:HOH:O	1.87	0.74
1:B:30:SER:HB3	1:B:174:LEU:CD2	2.18	0.73
1:F:181:GLN:O	1:F:186:TYR:HE1	1.70	0.73
1:F:73:GLN:HE21	1:F:135:LYS:H	1.34	0.73
1:C:165:GLU:HG3	3:C:962:HOH:O	1.89	0.73
1:A:168:ILE:HG22	1:A:172:ARG:NE	2.04	0.73
1:D:80:THR:OG1	1:D:83:GLU:HG3	1.88	0.73
1:F:163:SER:OG	1:F:165:GLU:HG2	1.88	0.73
1:B:73:GLN:HE21	1:B:135:LYS:H	1.36	0.72
1:A:37:VAL:HG13	1:A:159:SER:HB2	1.71	0.72
1:D:92:LEU:HD21	3:D:933:HOH:O	1.88	0.72
1:F:59:ASN:HB2	3:F:952:HOH:O	1.89	0.72
1:E:200:LEU:HD22	1:E:200:LEU:H	1.54	0.72
1:B:165:GLU:OE2	1:D:36:ARG:NH2	2.23	0.72
1:B:93:LEU:H	1:B:93:LEU:HD22	1.55	0.72
1:B:27:ILE:CD1	1:B:66:GLU:HG2	2.16	0.72
1:B:149:LEU:HD12	1:B:189:LEU:HD11	1.71	0.72
1:E:89:VAL:N	1:E:90:PRO:HD3	2.04	0.71
1:F:62:LYS:HB3	3:F:952:HOH:O	1.89	0.71
1:E:36:ARG:NH1	3:E:939:HOH:O	2.22	0.71
1:A:188:TYR:HE1	1:A:195:LYS:HZ1	1.38	0.71
1:C:133:TYR:HE1	3:C:992:HOH:O	1.73	0.70
1:E:174:LEU:HA	3:E:1032:HOH:O	1.89	0.70
1:A:181:GLN:CA	1:A:181:GLN:HE21	2.04	0.70
1:E:180:ILE:HG22	3:E:940:HOH:O	1.89	0.70
1:C:200:LEU:CD2	1:C:200:LEU:H	2.04	0.70
1:C:26:PRO:O	3:C:957:HOH:O	2.09	0.70
1:F:62:LYS:HE2	3:F:952:HOH:O	1.91	0.69
1:F:52:LYS:HE2	1:F:58:ARG:O	1.92	0.69
1:A:102:ILE:HD12	1:A:102:ILE:H	1.57	0.69
1:A:50:GLY:HA3	1:A:60:VAL:HG22	1.72	0.69
1:D:113:THR:HG23	1:D:184:LYS:HE3	1.74	0.69
1:A:93:LEU:H	1:A:93:LEU:HD22	1.56	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:190:HIS:O	1:B:191:GLU:CB	2.41	0.69
1:B:76:PHE:HB2	1:B:138:ILE:HG22	1.75	0.68
1:B:75:ILE:HG13	1:B:101:ILE:HG21	1.75	0.68
1:A:68:LYS:HB2	3:A:920:HOH:O	1.92	0.68
1:A:102:ILE:HD12	1:A:102:ILE:N	2.08	0.68
1:F:113:THR:HG23	3:F:971:HOH:O	1.93	0.68
1:D:30:SER:HB3	1:D:174:LEU:HD11	1.76	0.67
1:B:93:LEU:HD22	1:B:93:LEU:N	2.09	0.67
1:E:191:GLU:HB2	3:E:1054:HOH:O	1.94	0.67
1:D:116:ILE:O	1:D:120:CYS:HB2	1.94	0.67
1:E:191:GLU:HG3	1:E:195:LYS:CD	2.25	0.67
1:D:187:ASN:O	1:D:190:HIS:O	2.12	0.67
1:A:193:ARG:HA	3:A:944:HOH:O	1.95	0.67
1:B:193:ARG:HB2	3:B:995:HOH:O	1.94	0.67
1:A:37:VAL:HG12	1:A:37:VAL:O	1.94	0.67
1:A:168:ILE:O	1:A:172:ARG:HG3	1.96	0.66
1:F:160:ASP:N	1:F:160:ASP:OD2	2.28	0.66
1:C:188:TYR:HA	3:C:940:HOH:O	1.94	0.66
1:F:34:LEU:O	1:F:37:VAL:HG12	1.95	0.66
1:E:200:LEU:HD22	1:E:200:LEU:N	2.11	0.66
1:E:51:CYS:HA	3:E:980:HOH:O	1.96	0.66
1:F:116:ILE:CG2	1:F:195:LYS:HE2	2.26	0.66
1:A:27:ILE:HA	1:A:58:ARG:HH21	1.59	0.66
1:B:197:ALA:O	1:B:199:HIS:N	2.29	0.66
1:B:55:ASP:N	1:B:57:ARG:HH21	1.84	0.66
1:C:27:ILE:HG22	1:C:66:GLU:HG2	1.78	0.65
1:A:73:GLN:CB	1:A:134:ARG:HE	2.10	0.65
1:F:58:ARG:HH11	1:F:58:ARG:HG3	1.61	0.65
1:B:37:VAL:HG12	1:B:37:VAL:O	1.96	0.65
1:B:190:HIS:O	1:B:191:GLU:HB3	1.95	0.65
1:D:91:ASN:HD21	1:D:95:LEU:HD11	1.62	0.65
1:A:200:LEU:N	1:A:200:LEU:HD22	2.11	0.65
1:D:51:CYS:O	1:D:58:ARG:HG2	1.96	0.64
1:F:27:ILE:O	1:F:28:HIS:CG	2.50	0.64
1:F:162:ILE:HA	1:F:166:GLN:OE1	1.97	0.64
1:D:30:SER:CB	1:D:174:LEU:HD11	2.28	0.64
1:B:66:GLU:HG3	1:B:67:LEU:N	2.13	0.64
1:F:72:ILE:CD1	1:F:135:LYS:HE2	2.29	0.63
1:A:94:ASP:HB3	3:A:970:HOH:O	1.97	0.63
1:C:200:LEU:N	1:C:200:LEU:HD22	2.12	0.63
1:C:25:THR:HG23	1:C:26:PRO:HD2	1.81	0.63
1:E:197:ALA:O	1:E:199:HIS:N	2.31	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:60:VAL:HG21	1:E:89:VAL:HG12	1.81	0.63
1:F:73:GLN:HB3	1:F:134:ARG:NH1	2.12	0.63
1:D:83:GLU:HA	1:D:86:LYS:HG2	1.81	0.63
1:D:27:ILE:CG2	1:D:66:GLU:HG2	2.29	0.62
1:A:121:GLU:HG2	3:A:924:HOH:O	1.98	0.62
1:E:52:LYS:HD3	1:E:88:ARG:HH11	1.64	0.62
1:A:26:PRO:HB3	1:A:66:GLU:OE2	2.00	0.62
1:E:81:ARG:NH2	3:E:912:HOH:O	2.30	0.62
1:E:41:GLN:HE22	1:E:133:TYR:HA	1.64	0.62
1:F:80:THR:OG1	1:F:83:GLU:HG3	1.99	0.62
1:F:180:ILE:HG23	1:F:185:GLN:HB3	1.81	0.62
1:E:120:CYS:O	1:E:124:GLU:HG3	1.99	0.62
1:A:48:LEU:HD12	1:A:49:PRO:HD2	1.82	0.62
1:F:91:ASN:HD21	1:F:95:LEU:HD11	1.64	0.62
1:A:27:ILE:HG23	1:A:58:ARG:HE	1.65	0.62
1:C:116:ILE:HA	3:C:1021:HOH:O	1.99	0.62
1:D:174:LEU:HD12	1:D:175:ARG:HG2	1.81	0.62
1:F:27:ILE:C	1:F:28:HIS:CG	2.74	0.61
1:F:183:ILE:HD12	1:F:184:LYS:N	2.13	0.61
1:F:37:VAL:HG13	1:F:37:VAL:O	2.00	0.61
1:E:37:VAL:HG22	1:E:159:SER:HB2	1.82	0.61
1:E:52:LYS:HD3	1:E:88:ARG:NH1	2.15	0.61
1:F:179:ALA:HB2	3:F:969:HOH:O	2.01	0.61
1:C:33:SER:HB2	1:C:42:PHE:CE2	2.36	0.61
1:B:93:LEU:H	1:B:93:LEU:CD2	2.13	0.61
1:F:108:ILE:HD13	1:F:114:PRO:HD3	1.81	0.61
1:B:54:LYS:HA	1:B:57:ARG:NH2	2.16	0.61
1:D:27:ILE:HG22	1:D:66:GLU:HG2	1.81	0.61
1:C:149:LEU:HD13	3:C:960:HOH:O	2.00	0.61
1:A:73:GLN:NE2	1:A:134:ARG:HD2	2.16	0.61
1:F:60:VAL:CG1	1:F:89:VAL:HG12	2.31	0.60
1:F:179:ALA:O	1:F:181:GLN:N	2.33	0.60
1:E:73:GLN:HE21	1:E:135:LYS:H	1.49	0.60
1:D:37:VAL:O	1:D:37:VAL:HG13	2.00	0.60
1:C:50:GLY:HA3	1:C:60:VAL:HG13	1.82	0.60
1:B:55:ASP:N	1:B:57:ARG:NH2	2.44	0.60
1:F:148:CYS:HB3	3:F:926:HOH:O	2.02	0.60
1:D:46:CYS:O	3:D:1003:HOH:O	2.16	0.60
1:B:26:PRO:C	1:B:27:ILE:HD12	2.21	0.60
1:B:89:VAL:HG23	1:B:89:VAL:O	2.00	0.60
1:A:193:ARG:HH11	1:A:193:ARG:HG3	1.65	0.60
1:C:34:LEU:HB2	1:C:39:CYS:HB3	1.82	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:27:ILE:HD11	1:B:63:ASP:HA	1.84	0.60
1:F:133:TYR:HE1	3:F:972:HOH:O	1.83	0.59
1:B:62:LYS:HE2	3:B:941:HOH:O	2.01	0.59
1:F:91:ASN:HD22	1:F:95:LEU:CD2	2.11	0.59
1:A:79:CYS:HA	3:A:983:HOH:O	2.02	0.59
1:A:73:GLN:HB3	1:A:134:ARG:HH11	1.68	0.59
1:B:187:ASN:O	1:B:190:HIS:O	2.21	0.59
1:E:190:HIS:O	1:E:191:GLU:CB	2.51	0.59
1:E:159:SER:O	1:E:193:ARG:NH2	2.35	0.59
1:D:34:LEU:O	1:D:37:VAL:HG12	2.03	0.59
1:F:81:ARG:NH1	3:F:955:HOH:O	2.36	0.59
1:D:180:ILE:HG23	1:D:185:GLN:HB3	1.85	0.58
1:F:196:LEU:HD13	1:F:199:HIS:CE1	2.39	0.58
1:A:68:LYS:HE3	3:A:950:HOH:O	2.02	0.58
1:F:168:ILE:HG23	3:F:958:HOH:O	2.02	0.58
1:E:177:SER:HB3	3:E:931:HOH:O	2.03	0.58
1:F:192:PHE:C	1:F:192:PHE:CD1	2.76	0.58
1:E:191:GLU:CG	1:E:195:LYS:HD2	2.32	0.58
1:F:27:ILE:HD13	1:F:27:ILE:O	2.04	0.58
1:F:31:TRP:CH2	1:F:135:LYS:HE3	2.39	0.58
1:F:162:ILE:HG23	1:F:193:ARG:HH22	1.69	0.58
1:D:58:ARG:HG3	1:D:58:ARG:HH11	1.68	0.58
1:D:58:ARG:NH1	1:D:63:ASP:OD2	2.37	0.58
1:B:92:LEU:HD11	1:B:96:TYR:CZ	2.39	0.58
1:D:80:THR:N	3:D:923:HOH:O	2.36	0.58
1:D:66:GLU:HB3	3:D:922:HOH:O	2.02	0.58
1:B:93:LEU:HD12	1:B:103:THR:HG21	1.86	0.58
1:A:77:VAL:HG21	3:A:946:HOH:O	2.04	0.58
1:F:27:ILE:O	1:F:28:HIS:CD2	2.57	0.57
1:C:30:SER:OG	1:C:175:ARG:NH1	2.37	0.57
1:D:30:SER:HB3	1:D:174:LEU:CD1	2.34	0.57
1:E:118:SER:O	1:E:122:ILE:HG13	2.04	0.57
1:F:179:ALA:HB1	3:F:926:HOH:O	2.03	0.57
1:F:200:LEU:N	1:F:200:LEU:HD12	2.19	0.57
1:A:149:LEU:HG	3:A:954:HOH:O	2.04	0.57
1:A:148:CYS:SG	1:A:175:ARG:HG3	2.44	0.57
1:D:84:LEU:HD22	1:D:89:VAL:HG23	1.86	0.57
1:E:108:ILE:HD13	1:E:114:PRO:HD3	1.87	0.57
1:D:190:HIS:O	1:D:191:GLU:CG	2.52	0.57
1:F:81:ARG:HB2	1:F:81:ARG:CZ	2.34	0.57
1:A:164:PRO:O	1:A:168:ILE:HG13	2.05	0.57
1:A:188:TYR:HE1	1:A:195:LYS:NZ	2.01	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:81:ARG:HB2	1:F:81:ARG:HH11	1.70	0.56
1:A:50:GLY:HA2	1:A:58:ARG:HB3	1.86	0.56
1:F:56:VAL:C	1:F:57:ARG:HD3	2.25	0.56
1:A:102:ILE:HD13	1:A:134:ARG:NH1	2.20	0.56
1:A:197:ALA:O	1:A:199:HIS:N	2.39	0.56
1:C:172:ARG:HD3	1:C:177:SER:HA	1.86	0.56
1:A:144:LEU:HD11	1:A:178:GLY:C	2.26	0.56
1:F:112:GLY:HA2	3:F:971:HOH:O	2.05	0.56
1:D:58:ARG:NH2	3:D:970:HOH:O	2.39	0.56
1:C:84:LEU:CD1	1:C:93:LEU:HD21	2.35	0.55
1:A:85:SER:HA	1:A:90:PRO:HG3	1.88	0.55
1:E:36:ARG:NH1	1:E:36:ARG:HB2	2.11	0.55
1:E:36:ARG:HG2	3:E:993:HOH:O	2.07	0.55
1:F:70:CYS:O	1:F:135:LYS:NZ	2.28	0.55
1:F:190:HIS:O	1:F:191:GLU:HG2	2.06	0.55
1:C:187:ASN:O	1:C:190:HIS:O	2.24	0.55
1:A:94:ASP:O	1:A:98:GLN:HG3	2.06	0.55
1:E:36:ARG:HH11	1:E:36:ARG:CB	2.11	0.55
1:D:67:LEU:HG	3:D:922:HOH:O	2.06	0.55
1:F:149:LEU:N	1:F:180:ILE:HD11	2.21	0.55
1:B:149:LEU:HD12	1:B:189:LEU:CD1	2.36	0.55
1:C:133:TYR:N	1:C:133:TYR:CD1	2.74	0.55
1:D:58:ARG:CZ	1:D:58:ARG:HB3	2.37	0.55
1:A:86:LYS:HG2	1:A:87:TYR:CD2	2.41	0.55
1:A:67:LEU:O	1:A:72:ILE:HG12	2.06	0.54
1:F:36:ARG:HB2	3:F:945:HOH:O	2.06	0.54
1:F:181:GLN:CB	3:F:915:HOH:O	2.46	0.54
1:A:71:GLY:C	1:A:135:LYS:HD3	2.28	0.54
1:F:54:LYS:O	1:F:55:ASP:HB2	2.08	0.54
1:A:114:PRO:HG2	1:A:119:CYS:HB2	1.89	0.54
1:E:196:LEU:C	1:E:196:LEU:HD13	2.28	0.54
1:A:27:ILE:HD11	1:A:66:GLU:HG2	1.88	0.54
1:A:135:LYS:NZ	3:A:962:HOH:O	2.21	0.54
1:A:200:LEU:H	1:A:200:LEU:HD22	1.72	0.54
1:F:83:GLU:HA	1:F:86:LYS:CG	2.37	0.54
1:C:76:PHE:HB2	1:C:138:ILE:HG22	1.88	0.54
1:A:73:GLN:HG3	1:A:135:LYS:HB3	1.90	0.54
1:B:75:ILE:HG13	1:B:101:ILE:CG2	2.37	0.54
1:B:102:ILE:HD13	1:B:134:ARG:NH1	2.22	0.54
1:A:93:LEU:H	1:A:93:LEU:CD2	2.20	0.54
1:D:196:LEU:HD13	1:D:199:HIS:CE1	2.43	0.54
1:C:31:TRP:CG	3:C:988:HOH:O	2.60	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:163:SER:HB3	1:F:166:GLN:CG	2.33	0.54
1:A:193:ARG:NH1	1:A:193:ARG:HG3	2.23	0.54
1:C:131:LYS:HA	3:C:992:HOH:O	2.08	0.54
1:D:63:ASP:O	3:D:922:HOH:O	2.18	0.54
1:A:60:VAL:HG21	1:A:89:VAL:CG1	2.39	0.53
1:A:73:GLN:HB3	1:A:134:ARG:HE	1.73	0.53
1:F:26:PRO:O	1:F:27:ILE:HB	2.06	0.53
1:F:86:LYS:NZ	3:F:935:HOH:O	2.41	0.53
1:F:182:THR:O	1:F:186:TYR:HD1	1.92	0.53
1:D:58:ARG:NE	3:D:970:HOH:O	2.38	0.53
1:A:183:ILE:O	1:A:187:ASN:CG	2.46	0.53
1:F:73:GLN:HB3	1:F:134:ARG:CZ	2.38	0.53
1:F:102:ILE:CD1	1:F:102:ILE:N	2.72	0.53
1:F:60:VAL:HG23	1:F:61:GLN:N	2.23	0.53
1:A:75:ILE:HG12	1:A:137:LEU:HD23	1.90	0.53
1:A:102:ILE:CD1	1:A:134:ARG:HH12	2.21	0.53
1:A:187:ASN:O	1:A:190:HIS:O	2.27	0.53
1:E:190:HIS:O	1:E:191:GLU:HB3	2.08	0.53
1:D:121:GLU:O	1:D:125:GLU:HG3	2.08	0.53
1:A:149:LEU:HD13	1:A:185:GLN:HB3	1.88	0.53
1:E:190:HIS:O	1:E:191:GLU:HG2	2.08	0.53
1:B:37:VAL:O	1:B:37:VAL:CG1	2.57	0.53
1:D:68:LYS:HD3	1:D:99:CYS:O	2.09	0.53
1:F:156:LEU:HG	1:F:193:ARG:NH1	2.24	0.53
1:D:81:ARG:HB2	1:D:81:ARG:NH1	2.24	0.53
1:F:67:LEU:O	1:F:72:ILE:HG12	2.09	0.52
1:D:172:ARG:HD2	3:D:986:HOH:O	2.09	0.52
1:A:89:VAL:N	1:A:90:PRO:CD	2.72	0.52
1:E:172:ARG:HG3	3:E:1043:HOH:O	2.09	0.52
1:A:91:ASN:HB3	3:A:915:HOH:O	2.10	0.52
1:E:131:LYS:HE2	3:E:967:HOH:O	2.09	0.52
1:F:163:SER:N	1:F:166:GLN:OE1	2.36	0.52
1:C:137:LEU:HD12	1:C:138:ILE:N	2.25	0.52
1:E:33:SER:HB2	1:E:42:PHE:CE2	2.45	0.52
1:A:181:GLN:NE2	1:A:181:GLN:HA	2.10	0.52
1:C:50:GLY:HA3	1:C:60:VAL:CG1	2.40	0.52
1:B:54:LYS:O	1:B:55:ASP:HB2	2.09	0.52
1:F:156:LEU:HG	1:F:193:ARG:CZ	2.40	0.52
1:A:48:LEU:HD12	1:A:49:PRO:CD	2.39	0.52
1:F:116:ILE:CG2	1:F:195:LYS:HB3	2.40	0.51
1:E:33:SER:HB2	1:E:42:PHE:CZ	2.46	0.51
1:E:127:THR:CB	3:E:1059:HOH:O	2.46	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:113:THR:OG1	1:A:184:LYS:HE3	2.10	0.51
1:A:39:CYS:HB2	3:A:908:HOH:O	2.09	0.51
1:F:93:LEU:HD21	1:F:105:HIS:ND1	2.25	0.51
1:A:121:GLU:OE2	1:A:125:GLU:OE1	2.29	0.51
1:F:83:GLU:HA	1:F:86:LYS:HG2	1.93	0.51
1:C:148:CYS:C	1:C:180:ILE:HD11	2.31	0.51
1:F:163:SER:HG	1:F:165:GLU:HG2	1.76	0.51
1:C:37:VAL:O	1:C:37:VAL:CG1	2.58	0.51
1:F:91:ASN:O	1:F:95:LEU:HG	2.11	0.51
1:B:89:VAL:N	1:B:90:PRO:CD	2.74	0.51
1:A:196:LEU:HD13	1:A:196:LEU:O	2.11	0.51
1:E:165:GLU:O	1:E:169:ASP:OD1	2.29	0.51
1:D:52:LYS:HE3	3:D:936:HOH:O	2.11	0.51
1:C:73:GLN:NE2	3:C:925:HOH:O	2.44	0.51
1:B:91:ASN:O	1:B:95:LEU:HG	2.11	0.51
1:B:102:ILE:N	1:B:102:ILE:HD12	2.25	0.51
1:A:181:GLN:HB2	1:A:185:GLN:OE1	2.11	0.50
1:B:94:ASP:O	1:B:98:GLN:HG3	2.11	0.50
1:D:65:GLU:HG3	3:D:994:HOH:O	2.11	0.50
1:F:48:LEU:HD23	1:F:51:CYS:HB2	1.92	0.50
1:A:187:ASN:HB3	3:A:917:HOH:O	2.11	0.50
1:F:89:VAL:N	1:F:90:PRO:CD	2.72	0.50
1:E:60:VAL:CG2	1:E:89:VAL:HG12	2.41	0.50
1:F:86:LYS:HE2	1:F:87:TYR:CE2	2.47	0.50
1:A:180:ILE:HG21	1:A:186:TYR:CD2	2.47	0.50
1:F:23:GLU:OE1	1:F:23:GLU:N	2.45	0.50
1:A:190:HIS:O	1:A:191:GLU:CB	2.59	0.50
1:D:91:ASN:ND2	1:D:95:LEU:HG	2.27	0.50
1:F:172:ARG:NH2	3:F:958:HOH:O	2.45	0.49
1:A:33:SER:HB2	1:A:42:PHE:CE2	2.47	0.49
1:D:191:GLU:OE1	3:D:928:HOH:O	2.20	0.49
1:E:191:GLU:OE2	1:E:195:LYS:HD2	2.12	0.49
1:E:89:VAL:N	1:E:90:PRO:CD	2.75	0.49
1:A:93:LEU:HD22	1:A:93:LEU:N	2.23	0.49
1:E:149:LEU:HD12	1:E:185:GLN:HB3	1.94	0.49
1:F:81:ARG:HH11	1:F:81:ARG:CB	2.26	0.49
1:E:181:GLN:C	3:E:940:HOH:O	2.51	0.49
1:C:27:ILE:CG2	1:C:66:GLU:HG2	2.41	0.49
1:C:121:GLU:O	1:C:125:GLU:HG3	2.13	0.49
1:F:192:PHE:HB3	3:F:951:HOH:O	2.13	0.49
1:A:88:ARG:C	1:A:90:PRO:HD3	2.33	0.49
1:D:44:GLY:C	1:D:137:LEU:HD12	2.33	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:80:THR:O	1:F:84:LEU:HG	2.12	0.49
1:B:56:VAL:O	1:B:57:ARG:HD3	2.13	0.48
1:B:60:VAL:CG1	1:B:89:VAL:HG12	2.43	0.48
1:E:131:LYS:HD3	1:E:131:LYS:HA	1.49	0.48
1:A:146:ARG:CZ	3:A:983:HOH:O	2.61	0.48
1:F:175:ARG:HB2	3:F:969:HOH:O	2.13	0.48
1:A:146:ARG:NE	2:A:900:SO4:O1	2.36	0.48
1:D:41:GLN:HE22	1:D:133:TYR:HA	1.77	0.48
1:F:94:ASP:HA	1:F:97:GLN:HG3	1.95	0.48
1:D:89:VAL:N	1:D:90:PRO:HD3	2.28	0.48
1:A:139:HIS:ND1	1:A:139:HIS:C	2.67	0.48
1:C:84:LEU:HD13	1:C:93:LEU:HD21	1.94	0.48
1:E:88:ARG:C	1:E:90:PRO:HD3	2.34	0.48
1:F:58:ARG:NH1	1:F:58:ARG:HG3	2.25	0.48
1:E:41:GLN:NE2	1:E:133:TYR:HA	2.28	0.48
1:A:190:HIS:O	1:A:191:GLU:HB3	2.14	0.48
1:A:140:SER:HB3	1:A:141:TYR:H	1.53	0.48
1:E:182:THR:N	3:E:940:HOH:O	2.46	0.48
1:F:109:ALA:O	1:F:112:GLY:N	2.39	0.48
1:A:114:PRO:CG	1:A:119:CYS:HB2	2.44	0.48
1:A:25:THR:OG1	1:A:26:PRO:HD2	2.14	0.48
1:A:50:GLY:CA	1:A:60:VAL:HG22	2.42	0.48
1:A:193:ARG:CA	3:A:944:HOH:O	2.58	0.48
1:D:131:LYS:HD3	3:D:1009:HOH:O	2.13	0.48
1:F:164:PRO:N	3:F:999:HOH:O	2.46	0.48
1:B:60:VAL:HG11	1:B:89:VAL:HG12	1.96	0.48
1:E:37:VAL:O	1:E:37:VAL:CG1	2.62	0.48
1:D:143:GLY:CA	3:D:1003:HOH:O	2.62	0.48
1:F:178:GLY:O	1:F:179:ALA:O	2.31	0.47
1:B:37:VAL:HG11	1:B:158:LEU:HB2	1.96	0.47
1:C:144:LEU:N	3:C:968:HOH:O	2.46	0.47
1:C:60:VAL:HG22	3:C:913:HOH:O	2.14	0.47
1:E:115:ASP:OD1	1:E:117:ALA:N	2.47	0.47
1:E:187:ASN:O	1:E:190:HIS:O	2.32	0.47
1:D:58:ARG:CG	1:D:58:ARG:HH11	2.28	0.47
1:A:133:TYR:N	1:A:133:TYR:CD1	2.83	0.47
1:F:163:SER:CB	1:F:166:GLN:HG3	2.33	0.47
1:C:41:GLN:NE2	1:C:133:TYR:HA	2.29	0.47
1:B:181:GLN:HB3	1:B:185:GLN:OE1	2.15	0.47
1:E:26:PRO:O	1:E:58:ARG:NH2	2.34	0.47
1:F:156:LEU:HD11	3:F:999:HOH:O	2.14	0.47
1:F:93:LEU:O	1:F:97:GLN:HG2	2.13	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:91:ASN:ND2	1:D:95:LEU:CD1	2.78	0.47
1:D:26:PRO:C	1:D:27:ILE:HG22	2.35	0.47
1:C:34:LEU:CB	1:C:39:CYS:HB3	2.44	0.47
1:A:37:VAL:HG11	1:A:158:LEU:HB2	1.97	0.47
1:A:73:GLN:HE21	1:A:135:LYS:H	1.63	0.47
1:F:69:SER:O	3:F:931:HOH:O	2.20	0.47
1:D:195:LYS:HD2	3:D:954:HOH:O	2.15	0.47
1:A:156:LEU:HD21	1:A:189:LEU:HD12	1.96	0.47
1:B:119:CYS:O	1:B:123:MET:HG2	2.14	0.47
1:A:46:CYS:SG	1:A:137:LEU:HD11	2.53	0.47
1:C:165:GLU:OE2	1:C:190:HIS:NE2	2.48	0.47
1:A:73:GLN:HE21	1:A:134:ARG:HA	1.79	0.47
1:D:91:ASN:ND2	1:D:95:LEU:HD11	2.27	0.47
1:C:185:GLN:HB3	3:C:960:HOH:O	2.15	0.47
1:D:36:ARG:NE	3:D:927:HOH:O	2.48	0.47
1:D:144:LEU:HB2	3:D:951:HOH:O	2.15	0.47
1:F:182:THR:O	1:F:186:TYR:CD1	2.68	0.46
1:F:200:LEU:CD1	1:F:200:LEU:N	2.79	0.46
1:A:52:LYS:HA	1:A:56:VAL:O	2.15	0.46
1:B:116:ILE:HG21	1:B:195:LYS:HB3	1.96	0.46
1:B:88:ARG:C	1:B:90:PRO:HD3	2.35	0.46
1:F:163:SER:H	1:F:166:GLN:CD	2.17	0.46
1:E:50:GLY:HA3	1:E:60:VAL:HG22	1.98	0.46
1:D:47:ALA:O	1:D:49:PRO:HD3	2.16	0.46
1:F:27:ILE:C	1:F:28:HIS:ND1	2.69	0.46
1:F:81:ARG:NH1	1:F:81:ARG:CB	2.73	0.46
1:D:91:ASN:HD21	1:D:95:LEU:CD1	2.29	0.46
1:D:91:ASN:HD22	1:D:95:LEU:HG	1.80	0.46
1:F:83:GLU:N	3:F:977:HOH:O	2.48	0.46
1:C:25:THR:CG2	1:C:26:PRO:HD2	2.45	0.46
1:A:102:ILE:HD11	1:A:134:ARG:HH12	1.81	0.46
1:A:193:ARG:CB	3:A:903:HOH:O	2.63	0.46
1:C:100:GLY:HA2	3:C:987:HOH:O	2.16	0.46
1:E:184:LYS:HG3	3:E:964:HOH:O	2.15	0.46
1:A:60:VAL:HG21	1:A:89:VAL:HG12	1.98	0.46
1:E:198:ALA:HB3	3:E:937:HOH:O	2.15	0.46
1:C:91:ASN:HB2	3:C:956:HOH:O	2.15	0.46
1:F:60:VAL:HG12	1:F:89:VAL:HG12	1.98	0.46
1:D:190:HIS:O	1:D:191:GLU:CB	2.63	0.46
1:C:193:ARG:HD2	3:C:934:HOH:O	2.15	0.46
1:B:190:HIS:O	1:B:191:GLU:HG2	2.16	0.46
1:A:102:ILE:CD1	1:A:134:ARG:NH1	2.79	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:174:LEU:C	1:D:174:LEU:HD13	2.36	0.46
1:E:196:LEU:O	1:E:196:LEU:HD13	2.16	0.46
1:C:189:LEU:HD12	1:C:189:LEU:HA	1.82	0.46
1:F:181:GLN:O	1:F:186:TYR:CE1	2.59	0.45
1:A:31:TRP:CH2	1:A:135:LYS:HE3	2.50	0.45
1:C:31:TRP:CD1	3:C:988:HOH:O	2.56	0.45
1:F:181:GLN:OE1	1:F:185:GLN:OE1	2.34	0.45
1:F:163:SER:C	3:F:999:HOH:O	2.54	0.45
1:A:72:ILE:HD13	1:A:135:LYS:HE2	1.97	0.45
1:E:58:ARG:HH21	1:E:63:ASP:CG	2.20	0.45
1:D:178:GLY:O	1:D:179:ALA:O	2.33	0.45
1:A:149:LEU:HD12	1:A:189:LEU:HD22	1.99	0.45
1:E:36:ARG:HB2	3:E:1024:HOH:O	2.17	0.45
1:B:190:HIS:O	1:B:191:GLU:CG	2.65	0.45
1:D:180:ILE:H	1:D:180:ILE:HD12	1.81	0.45
1:A:78:PHE:CZ	1:A:150:VAL:HG21	2.52	0.45
1:A:37:VAL:HB	3:A:908:HOH:O	2.17	0.45
1:E:81:ARG:HG2	3:E:921:HOH:O	2.16	0.45
1:E:156:LEU:HD11	1:E:167:ALA:CB	2.47	0.45
1:E:37:VAL:O	1:E:37:VAL:HG13	2.17	0.45
1:E:104:HIS:N	1:E:104:HIS:CD2	2.85	0.45
1:D:154:LEU:O	1:D:157:TYR:HB3	2.17	0.45
1:D:93:LEU:N	1:D:93:LEU:CD1	2.80	0.45
1:A:116:ILE:O	1:A:120:CYS:HB2	2.17	0.45
1:D:199:HIS:HB2	3:D:1011:HOH:O	2.17	0.45
1:A:153:CYS:SG	3:A:954:HOH:O	2.62	0.44
1:E:200:LEU:H	1:E:200:LEU:CD2	2.27	0.44
1:E:85:SER:HA	1:E:90:PRO:HG3	1.99	0.44
1:A:164:PRO:HD3	3:A:943:HOH:O	2.17	0.44
1:D:27:ILE:HG23	1:D:28:HIS:N	2.33	0.44
1:A:144:LEU:HD12	1:A:179:ALA:HA	1.98	0.44
1:F:36:ARG:CB	3:F:945:HOH:O	2.63	0.44
1:D:48:LEU:HD12	3:D:933:HOH:O	2.16	0.44
1:F:149:LEU:HD12	1:F:149:LEU:O	2.17	0.44
1:E:190:HIS:O	1:E:191:GLU:CG	2.65	0.44
1:C:52:LYS:NZ	3:C:913:HOH:O	2.49	0.44
1:D:99:CYS:HA	3:D:946:HOH:O	2.16	0.44
1:C:196:LEU:O	1:C:197:ALA:C	2.56	0.44
1:A:25:THR:OG1	1:A:26:PRO:CD	2.66	0.44
1:B:165:GLU:HB2	3:B:905:HOH:O	2.17	0.44
1:A:193:ARG:HD3	3:A:944:HOH:O	2.18	0.44
1:A:72:ILE:HA	1:A:135:LYS:HG2	1.98	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:102:ILE:HD12	1:F:102:ILE:N	2.33	0.44
1:B:163:SER:OG	1:B:166:GLN:HG3	2.18	0.44
1:B:27:ILE:HA	1:B:58:ARG:HH22	1.82	0.44
1:A:82:GLY:O	1:A:85:SER:HB2	2.17	0.44
1:A:149:LEU:HD12	1:A:189:LEU:CD2	2.47	0.44
1:B:116:ILE:HD13	3:B:933:HOH:O	2.16	0.44
1:C:89:VAL:HG23	1:C:89:VAL:O	2.18	0.44
1:F:29:ILE:HG21	1:F:31:TRP:CZ2	2.53	0.44
1:A:33:SER:HB2	1:A:42:PHE:CZ	2.53	0.44
1:F:63:ASP:O	1:F:66:GLU:HB3	2.18	0.43
1:D:27:ILE:HG21	1:D:66:GLU:HG2	2.00	0.43
1:F:187:ASN:O	1:F:190:HIS:O	2.36	0.43
1:D:82:GLY:HA3	3:D:984:HOH:O	2.18	0.43
1:A:54:LYS:HB2	1:A:54:LYS:HE3	1.80	0.43
1:F:131:LYS:HG3	1:F:132:ASN:OD1	2.18	0.43
1:F:60:VAL:HG11	1:F:89:VAL:HG12	2.00	0.43
1:B:85:SER:HA	1:B:90:PRO:HG3	2.01	0.43
1:D:27:ILE:HG12	3:D:977:HOH:O	2.18	0.43
1:C:36:ARG:NH1	3:C:911:HOH:O	2.50	0.43
1:F:47:ALA:HB1	3:F:989:HOH:O	2.18	0.43
1:F:94:ASP:CB	3:F:978:HOH:O	2.53	0.43
1:A:73:GLN:HB2	1:A:134:ARG:HE	1.83	0.43
1:C:55:ASP:HB3	3:C:1034:HOH:O	2.17	0.43
1:F:91:ASN:ND2	1:F:95:LEU:HD21	2.19	0.43
1:C:116:ILE:HG23	3:C:1021:HOH:O	2.17	0.43
1:A:165:GLU:OE1	1:A:190:HIS:NE2	2.51	0.43
1:A:163:SER:OG	1:A:166:GLN:HG3	2.18	0.43
1:E:102:ILE:N	1:E:102:ILE:HD12	2.34	0.43
1:C:84:LEU:HD11	1:C:93:LEU:HD21	1.99	0.43
1:F:116:ILE:HG21	1:F:195:LYS:HB3	2.01	0.43
1:E:200:LEU:N	1:E:200:LEU:CD2	2.82	0.43
1:A:102:ILE:N	1:A:102:ILE:CD1	2.79	0.43
1:C:38:ASN:HA	1:C:38:ASN:HD22	1.64	0.43
1:F:144:LEU:HD12	1:F:179:ALA:HA	2.00	0.43
1:D:30:SER:HB2	1:D:174:LEU:HD11	1.99	0.43
1:F:37:VAL:HB	1:F:155:LEU:HD22	2.01	0.43
1:B:195:LYS:O	1:B:198:ALA:HB3	2.19	0.42
1:F:196:LEU:CD1	1:F:199:HIS:CE1	3.02	0.42
1:A:78:PHE:CE2	1:A:150:VAL:HG21	2.54	0.42
1:B:50:GLY:HA3	1:B:60:VAL:CG1	2.49	0.42
1:C:144:LEU:HB3	3:C:923:HOH:O	2.18	0.42
1:F:68:LYS:HA	1:F:68:LYS:HD2	1.80	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:92:LEU:HD11	1:F:96:TYR:OH	2.20	0.42
1:A:163:SER:HA	3:A:943:HOH:O	2.19	0.42
1:C:133:TYR:N	1:C:133:TYR:HD1	2.15	0.42
1:C:37:VAL:HG11	1:C:158:LEU:HB2	2.01	0.42
1:F:29:ILE:HD13	1:F:137:LEU:HD13	2.02	0.42
1:C:82:GLY:N	3:C:937:HOH:O	2.44	0.42
1:F:186:TYR:HE2	3:F:906:HOH:O	2.03	0.42
1:F:153:CYS:HG	1:F:192:PHE:HD2	1.65	0.42
1:C:178:GLY:O	1:C:179:ALA:O	2.37	0.42
1:A:37:VAL:CG1	1:A:37:VAL:O	2.67	0.42
1:F:66:GLU:HB2	3:F:970:HOH:O	2.19	0.42
1:D:86:LYS:HZ3	1:D:86:LYS:HB2	1.85	0.42
1:A:193:ARG:HB2	3:A:903:HOH:O	2.20	0.42
1:F:198:ALA:HA	1:F:200:LEU:CD1	2.50	0.42
1:B:200:LEU:HD22	1:B:200:LEU:N	2.35	0.42
1:B:27:ILE:HD13	1:B:66:GLU:CG	2.23	0.42
1:A:27:ILE:HA	1:A:58:ARG:NH2	2.30	0.42
1:C:196:LEU:HD22	1:C:196:LEU:HA	1.76	0.42
1:E:93:LEU:HD21	1:E:105:HIS:ND1	2.35	0.42
1:F:61:GLN:H	1:F:61:GLN:HG2	1.27	0.41
1:F:162:ILE:HG12	1:F:163:SER:N	2.34	0.41
1:F:148:CYS:C	1:F:180:ILE:HD11	2.40	0.41
1:C:190:HIS:O	1:C:191:GLU:CG	2.62	0.41
1:E:180:ILE:HG12	3:E:1043:HOH:O	2.20	0.41
1:C:131:LYS:HD3	1:C:131:LYS:HA	1.79	0.41
1:A:134:ARG:NH2	3:A:988:HOH:O	2.53	0.41
1:B:91:ASN:OD1	1:B:91:ASN:N	2.51	0.41
1:B:131:LYS:HA	1:B:131:LYS:HD3	1.69	0.41
1:A:60:VAL:CG2	1:A:89:VAL:CG1	2.98	0.41
1:A:53:PHE:CE2	1:A:54:LYS:HE2	2.55	0.41
1:A:189:LEU:HA	1:A:189:LEU:HD13	1.91	0.41
1:F:192:PHE:HD1	1:F:193:ARG:N	2.18	0.41
1:F:112:GLY:C	3:F:971:HOH:O	2.59	0.41
1:A:30:SER:OG	1:A:175:ARG:NH1	2.53	0.41
1:A:180:ILE:HB	1:A:186:TYR:CE2	2.55	0.41
1:F:172:ARG:CZ	3:F:958:HOH:O	2.68	0.41
1:F:180:ILE:HB	3:F:958:HOH:O	2.20	0.41
1:C:102:ILE:HD13	3:C:994:HOH:O	2.21	0.41
1:C:29:ILE:C	3:C:988:HOH:O	2.59	0.41
1:F:63:ASP:O	1:F:67:LEU:HG	2.20	0.41
1:A:200:LEU:CD2	1:A:200:LEU:N	2.82	0.41
1:F:91:ASN:HD22	1:F:95:LEU:CG	2.34	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:191:GLU:OE2	1:B:195:LYS:HE3	2.21	0.41
1:A:169:ASP:HA	1:A:172:ARG:HD2	2.03	0.41
1:C:33:SER:HA	1:C:42:PHE:CD2	2.56	0.41
1:B:92:LEU:HD11	1:B:96:TYR:CE2	2.55	0.41
1:B:102:ILE:HD13	1:B:134:ARG:HH11	1.86	0.41
1:A:190:HIS:O	1:A:191:GLU:HG2	2.21	0.41
1:E:184:LYS:HB3	1:E:184:LYS:HE3	1.79	0.41
1:A:75:ILE:CD1	1:A:101:ILE:HG21	2.51	0.41
1:A:168:ILE:HG22	1:A:172:ARG:HE	1.81	0.41
1:C:41:GLN:HG3	3:C:1035:HOH:O	2.21	0.41
1:F:183:ILE:HD12	1:F:183:ILE:C	2.41	0.41
1:F:184:LYS:HD3	3:F:971:HOH:O	2.21	0.41
1:C:118:SER:O	1:C:122:ILE:HG13	2.20	0.41
1:F:44:GLY:C	1:F:45:LEU:HD12	2.41	0.41
1:D:174:LEU:O	1:D:174:LEU:HD13	2.21	0.40
1:B:50:GLY:HA3	1:B:60:VAL:HG12	2.03	0.40
1:A:49:PRO:HG2	1:A:96:TYR:OH	2.21	0.40
1:F:196:LEU:O	1:F:197:ALA:C	2.60	0.40
1:D:81:ARG:NH2	3:D:930:HOH:O	2.54	0.40
1:D:179:ALA:O	1:D:181:GLN:N	2.54	0.40
1:D:54:LYS:O	1:D:55:ASP:HB2	2.22	0.40
1:C:26:PRO:C	1:C:27:ILE:HG22	2.42	0.40
1:D:196:LEU:O	1:D:197:ALA:C	2.59	0.40
1:A:91:ASN:HB3	3:A:937:HOH:O	2.21	0.40
1:E:119:CYS:O	1:E:123:MET:HG2	2.21	0.40
1:E:168:ILE:CG2	3:E:1043:HOH:O	2.59	0.40
1:D:196:LEU:O	1:D:199:HIS:CE1	2.75	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:24:GLN:OE1	3:E:931:HOH:O[5_565]	2.06	0.14
1:A:61:GLN:OE1	1:A:181:GLN:O[6_554]	2.14	0.06

## 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	174/212 (82%)	159 (91%)	11 (6%)	4 (2%)	10	3
1	B	174/212 (82%)	165 (95%)	6 (3%)	3 (2%)	14	5
1	C	176/212 (83%)	164 (93%)	5 (3%)	7 (4%)	5	1
1	D	176/212 (83%)	158 (90%)	11 (6%)	7 (4%)	5	1
1	E	174/212 (82%)	165 (95%)	5 (3%)	4 (2%)	10	3
1	F	176/212 (83%)	154 (88%)	15 (8%)	7 (4%)	5	1
All	All	1050/1272 (82%)	965 (92%)	53 (5%)	32 (3%)	7	2

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	26	PRO
1	B	26	PRO
1	B	198	ALA
1	C	27	ILE
1	C	54	LYS
1	C	179	ALA
1	D	27	ILE
1	D	179	ALA
1	F	27	ILE
1	F	54	LYS
1	F	179	ALA
1	A	198	ALA
1	B	191	GLU
1	C	28	HIS
1	D	54	LYS
1	E	26	PRO
1	E	198	ALA
1	F	144	LEU
1	F	180	ILE
1	A	191	GLU
1	C	26	PRO
1	D	26	PRO
1	F	28	HIS
1	C	144	LEU
1	D	28	HIS
1	D	144	LEU
1	E	27	ILE
1	E	191	GLU

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Mol	Chain	Res	Type
1	A	27	ILE
1	C	180	ILE
1	D	180	ILE
1	F	26	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	155/191 (81%)	133 (86%)	22 (14%)	5	2
1	B	155/191 (81%)	146 (94%)	9 (6%)	28	21
1	C	157/191 (82%)	139 (88%)	18 (12%)	8	4
1	D	157/191 (82%)	147 (94%)	10 (6%)	25	17
1	E	155/191 (81%)	142 (92%)	13 (8%)	16	9
1	F	157/191 (82%)	136 (87%)	21 (13%)	6	3
All	All	936/1146 (82%)	843 (90%)	93 (10%)	11	6

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ILE
1	A	55	ASP
1	A	58	ARG
1	A	63	ASP
1	A	81	ARG
1	A	89	VAL
1	A	99	CYS
1	A	121	GLU
1	A	133	TYR
1	A	137	LEU
1	A	139	HIS
1	A	140	SER
1	A	156	LEU
1	A	174	LEU
1	A	181	GLN

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Mol	Chain	Res	Type
1	A	182	THR
1	A	186	TYR
1	A	189	LEU
1	A	192	PHE
1	A	193	ARG
1	A	195	LYS
1	A	200	LEU
1	B	55	ASP
1	B	66	GLU
1	B	81	ARG
1	B	85	SER
1	B	99	CYS
1	B	156	LEU
1	B	174	LEU
1	B	189	LEU
1	B	200	LEU
1	C	23	GLU
1	C	26	PRO
1	C	33	SER
1	C	36	ARG
1	C	37	VAL
1	C	38	ASN
1	C	45	LEU
1	C	52	LYS
1	C	54	LYS
1	C	60	VAL
1	C	93	LEU
1	C	140	SER
1	C	147	SER
1	C	156	LEU
1	C	174	LEU
1	C	189	LEU
1	C	196	LEU
1	C	199	HIS
1	D	26	PRO
1	D	37	VAL
1	D	58	ARG
1	D	65	GLU
1	D	81	ARG
1	D	116	ILE
1	D	147	SER
1	D	156	LEU

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Mol	Chain	Res	Type
1	D	165	GLU
1	D	189	LEU
1	E	36	ARG
1	E	38	ASN
1	E	58	ARG
1	E	66	GLU
1	E	70	CYS
1	E	93	LEU
1	E	99	CYS
1	E	121	GLU
1	E	140	SER
1	E	161	THR
1	E	174	LEU
1	E	189	LEU
1	E	200	LEU
1	F	23	GLU
1	F	24	GLN
1	F	26	PRO
1	F	27	ILE
1	F	28	HIS
1	F	30	SER
1	F	37	VAL
1	F	61	GLN
1	F	63	ASP
1	F	81	ARG
1	F	97	GLN
1	F	102	ILE
1	F	138	ILE
1	F	149	LEU
1	F	160	ASP
1	F	162	ILE
1	F	165	GLU
1	F	174	LEU
1	F	181	GLN
1	F	192	PHE
1	F	200	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	98	GLN

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Mol	Chain	Res	Type
1	A	181	GLN
1	A	187	ASN
1	B	38	ASN
1	B	41	GLN
1	B	73	GLN
1	B	98	GLN
1	B	187	ASN
1	B	190	HIS
1	C	24	GLN
1	C	38	ASN
1	C	41	GLN
1	C	73	GLN
1	C	187	ASN
1	D	28	HIS
1	D	38	ASN
1	D	41	GLN
1	D	73	GLN
1	D	91	ASN
1	D	97	GLN
1	D	187	ASN
1	E	38	ASN
1	E	41	GLN
1	E	73	GLN
1	E	187	ASN
1	F	73	GLN
1	F	91	ASN
1	F	97	GLN
1	F	187	ASN
1	F	199	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

6 ligands 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	$\# Z  > 2$
2	SO4	A	900	-	4,4,4	0.35	0	6,6,6	0.49	0
2	SO4	B	901	-	4,4,4	0.49	0	6,6,6	0.23	0
2	SO4	C	902	-	4,4,4	0.56	0	6,6,6	0.25	0
2	SO4	D	903	-	4,4,4	0.32	0	6,6,6	0.32	0
2	SO4	E	904	-	4,4,4	0.15	0	6,6,6	0.28	0
2	SO4	F	905	-	4,4,4	0.26	0	6,6,6	0.14	0

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
2	SO4	A	900	-	-	0/0/0/0	0/0/0/0
2	SO4	B	901	-	-	0/0/0/0	0/0/0/0
2	SO4	C	902	-	-	0/0/0/0	0/0/0/0
2	SO4	D	903	-	-	0/0/0/0	0/0/0/0
2	SO4	E	904	-	-	0/0/0/0	0/0/0/0
2	SO4	F	905	-	-	0/0/0/0	0/0/0/0

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.

## 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	176/212 (83%)	0.63	15 (8%) 11 10	14, 44, 75, 78	0
1	B	176/212 (83%)	-0.28	2 (1%) 77 78	10, 21, 45, 64	0
1	C	178/212 (83%)	-0.21	0 100 100	13, 23, 41, 58	0
1	D	178/212 (83%)	-0.18	1 (0%) 86 88	10, 25, 42, 63	0
1	E	176/212 (83%)	-0.27	1 (0%) 86 88	10, 21, 42, 59	0
1	F	178/212 (83%)	0.63	17 (9%) 8 8	13, 45, 74, 83	0
All	All	1062/1272 (83%)	0.05	36 (3%) 43 43	10, 26, 63, 83	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	183	ILE	5.3
1	F	198	ALA	5.1
1	F	183	ILE	4.9
1	A	56	VAL	4.6
1	A	178	GLY	4.6
1	F	57	ARG	4.1
1	F	56	VAL	3.5
1	A	143	GLY	3.5
1	F	181	GLN	3.5
1	A	182	THR	3.3
1	A	176	GLY	3.3
1	A	111	GLY	3.2
1	F	50	GLY	3.1
1	A	179	ALA	3.0
1	A	184	LYS	3.0
1	D	57	ARG	2.9
1	F	54	LYS	2.9
1	A	180	ILE	2.8
1	F	27	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	89	VAL	2.4
1	F	89	VAL	2.4
1	F	59	ASN	2.4
1	F	176	GLY	2.4
1	A	192	PHE	2.4
1	A	60	VAL	2.3
1	F	25	THR	2.3
1	E	55	ASP	2.3
1	F	55	ASP	2.3
1	F	37	VAL	2.2
1	F	49	PRO	2.2
1	B	55	ASP	2.1
1	A	81	ARG	2.1
1	F	191	GLU	2.1
1	F	60	VAL	2.1
1	A	187	ASN	2.0
1	B	56	VAL	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 ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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
2	SO4	E	904	5/5	0.12	1.02	30,32,33,34	0
2	SO4	B	901	5/5	0.11	0.16	11,12,14,16	0
2	SO4	D	903	5/5	0.11	-0.32	15,15,15,18	0
2	SO4	A	900	5/5	0.15	-0.41	28,32,33,34	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	C	902	5/5	0.10	-0.45	17,17,18,21	0
2	SO4	F	905	5/5	0.12	-0.88	55,56,57,57	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.