



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 01:09 AM GMT

PDB ID : 2OCP
Title : Crystal Structure of Human Deoxyguanosine Kinase
Authors : Johansson, K.; Ramaswamy, S.; Ljungkrantz, C.; Knecht, W.; Piskur, J.;
Munch-Petersen, B.; Eriksson, S.; Eklund, H.
Deposited on : 2006-12-21
Resolution : 2.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

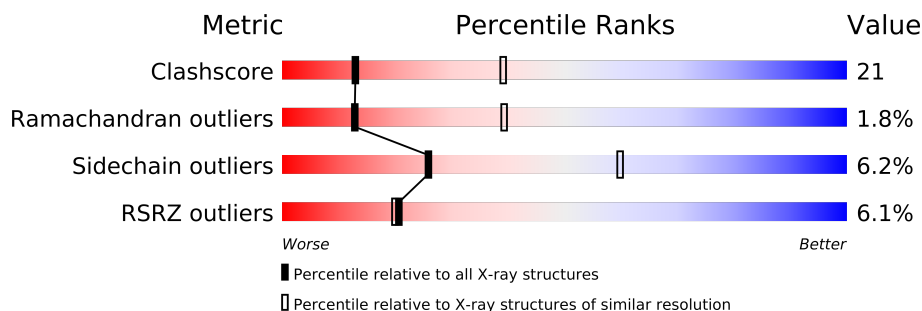
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.80 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))
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	241	
1	B	241	
1	C	241	
1	D	241	
1	E	241	
1	F	241	
1	G	241	
1	H	241	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15832 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.

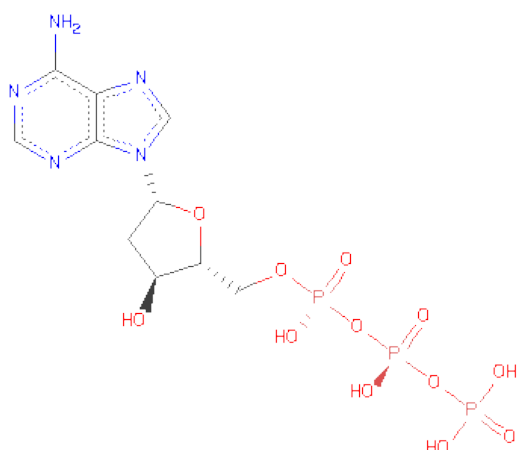
- Molecule 1 is a protein called Deoxyguanosine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			
1	B	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			
1	C	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			
1	D	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			
1	E	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			
1	F	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			
1	G	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			
1	H	229	Total	C	N	O	S	0	0	0
			1924	1249	326	344	5			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	83	ASN	THR	variant	UNP Q16854
B	83	ASN	THR	variant	UNP Q16854
C	83	ASN	THR	variant	UNP Q16854
D	83	ASN	THR	variant	UNP Q16854
E	83	ASN	THR	variant	UNP Q16854
F	83	ASN	THR	variant	UNP Q16854
G	83	ASN	THR	variant	UNP Q16854
H	83	ASN	THR	variant	UNP Q16854

- Molecule 2 is 2'-DEOXYADENOSINE 5'-TRIPHOSPHATE (three-letter code: DTP) (formula: $C_{10}H_{16}N_5O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
2	B	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
2	C	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
2	D	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
2	E	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
2	F	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
2	G	1	Total	C	N	O	P	0	0
			30	10	5	12	3		
2	H	1	Total	C	N	O	P	0	0
			30	10	5	12	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	25	Total	O	0	0
			25	25		
3	B	25	Total	O	0	0
			25	25		
3	C	25	Total	O	0	0
			25	25		
3	D	25	Total	O	0	0
			25	25		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	25	Total 25	O 25	0	0
3	F	25	Total 25	O 25	0	0
3	G	25	Total 25	O 25	0	0
3	H	25	Total 25	O 25	0	0

3 Residue-property plots

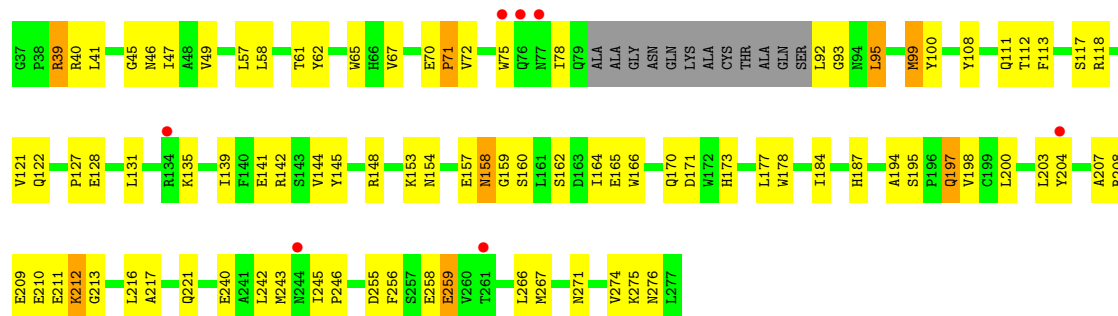
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Deoxyguanosine kinase



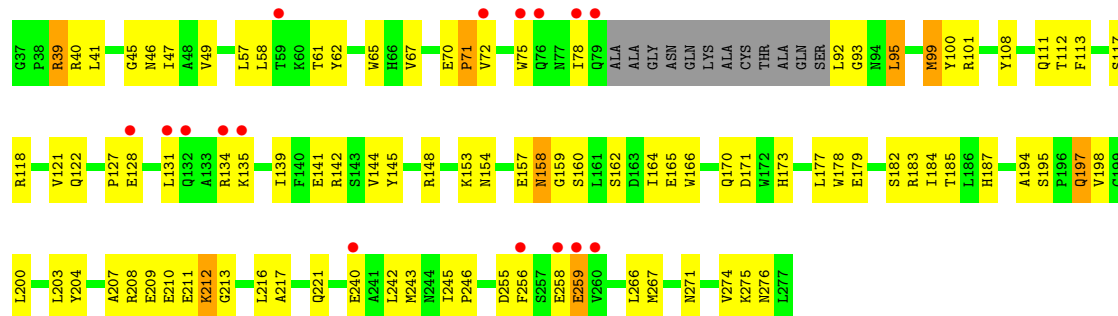
- Molecule 1: Deoxyguanosine kinase

Chain D:



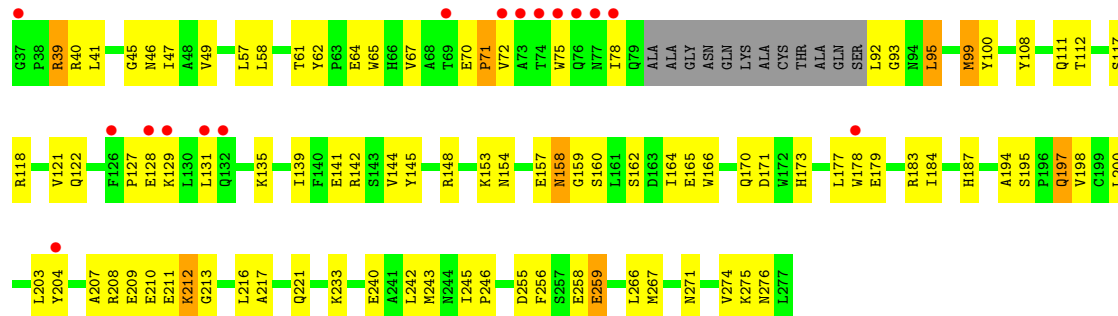
- Molecule 1: Deoxyguanosine kinase

Chain E:



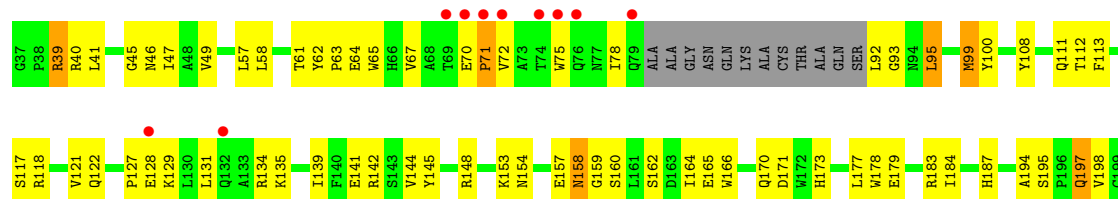
- Molecule 1: Deoxyguanosine kinase

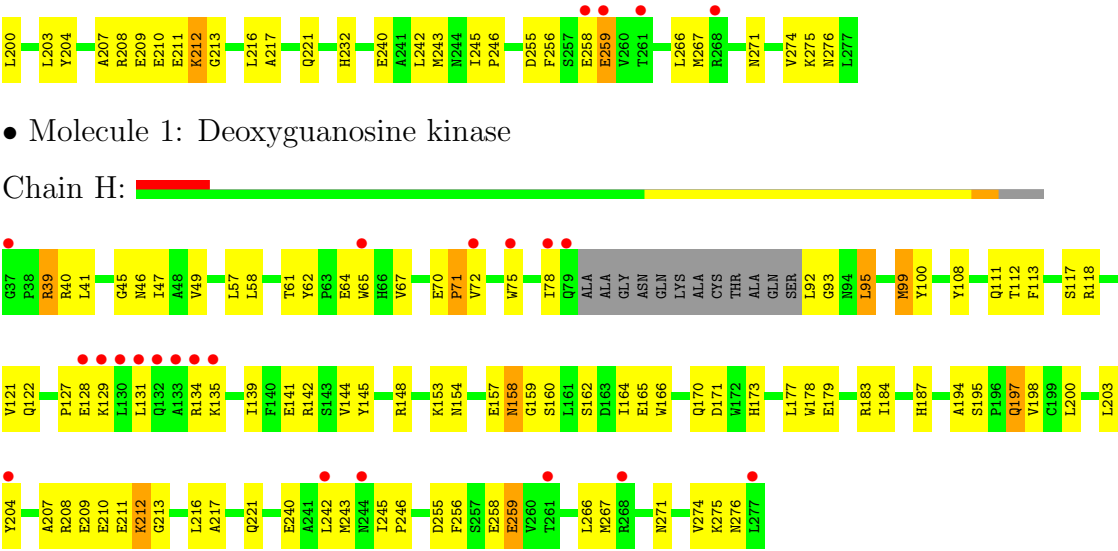
Chain F:



- Molecule 1: Deoxyguanosine kinase

Chain G:





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.18Å 120.03Å 154.35Å 90.00° 90.68° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 38.77 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.5 (30.00-2.80) 97.5 (38.77-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.81Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.265 , 0.277 0.285 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	35.1	Xtriage
Anisotropy	0.976	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 10.8	EDS
Estimated twinning fraction	0.035 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 59766 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	15832	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.84% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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: DTP

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	0.40	0/1976	0.61	0/2676
1	B	0.40	0/1976	0.61	0/2676
1	C	0.40	0/1976	0.61	0/2676
1	D	0.40	0/1976	0.61	0/2676
1	E	0.40	0/1976	0.61	0/2676
1	F	0.40	0/1976	0.61	0/2676
1	G	0.40	0/1976	0.61	0/2676
1	H	0.40	0/1976	0.61	0/2676
All	All	0.40	0/15808	0.61	0/21408

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1924	0	1899	81	1
1	B	1924	0	1899	90	3
1	C	1924	0	1899	76	10
1	D	1924	0	1899	77	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1924	0	1899	87	11
1	F	1924	0	1899	83	3
1	G	1924	0	1899	83	7
1	H	1924	0	1899	82	3
2	A	30	0	12	1	0
2	B	30	0	12	3	0
2	C	30	0	12	2	0
2	D	30	0	12	2	0
2	E	30	0	12	1	0
2	F	30	0	12	2	0
2	G	30	0	12	3	0
2	H	30	0	12	3	0
3	A	25	0	0	5	0
3	B	25	0	0	5	0
3	C	25	0	0	5	0
3	D	25	0	0	5	0
3	E	25	0	0	5	0
3	F	25	0	0	5	0
3	G	25	0	0	5	0
3	H	25	0	0	5	0
All	All	15832	0	15288	639	20

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 21.

All (639) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:244:ASN:CB	1:E:134:ARG:HH22	1.36	1.34
1:B:244:ASN:HB3	1:E:134:ARG:NH2	1.58	1.16
1:B:244:ASN:CB	1:E:134:ARG:NH2	2.13	1.08
1:H:40:ARG:H	1:H:187:HIS:HD2	1.01	1.00
1:G:40:ARG:H	1:G:187:HIS:HD2	1.01	1.00
1:B:244:ASN:HB3	1:E:134:ARG:HH22	0.84	0.99
1:E:40:ARG:H	1:E:187:HIS:HD2	1.01	0.99
1:B:40:ARG:H	1:B:187:HIS:HD2	1.01	0.98
1:C:40:ARG:H	1:C:187:HIS:HD2	1.01	0.95
1:A:40:ARG:H	1:A:187:HIS:HD2	1.01	0.93
1:F:40:ARG:H	1:F:187:HIS:HD2	1.01	0.93
1:D:40:ARG:H	1:D:187:HIS:HD2	1.01	0.92
1:C:40:ARG:H	1:C:187:HIS:CD2	1.88	0.92
1:A:40:ARG:H	1:A:187:HIS:CD2	1.88	0.92
1:E:40:ARG:H	1:E:187:HIS:CD2	1.88	0.91

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:40:ARG:H	1:D:187:HIS:CD2	1.88	0.91
1:B:40:ARG:H	1:B:187:HIS:CD2	1.88	0.91
1:F:40:ARG:H	1:F:187:HIS:CD2	1.88	0.91
1:G:40:ARG:H	1:G:187:HIS:CD2	1.88	0.90
1:B:244:ASN:CG	1:E:134:ARG:HH22	1.75	0.90
1:H:40:ARG:H	1:H:187:HIS:CD2	1.88	0.90
1:B:244:ASN:CG	1:E:134:ARG:NH2	2.28	0.86
1:C:58:LEU:HD23	1:C:267:MET:HE1	1.57	0.85
1:F:118:ARG:HH12	2:F:301:DTP:HN61	1.25	0.84
1:A:118:ARG:HH12	2:A:301:DTP:HN61	1.26	0.83
1:G:118:ARG:HH12	2:G:301:DTP:HN61	1.27	0.83
1:B:118:ARG:HH12	2:B:301:DTP:HN61	1.27	0.82
1:H:58:LEU:HD23	1:H:267:MET:HE1	1.62	0.81
1:D:118:ARG:HH12	2:D:301:DTP:HN61	1.23	0.81
1:F:58:LEU:HD23	1:F:267:MET:HE1	1.62	0.81
1:E:118:ARG:HH12	2:E:301:DTP:HN61	1.28	0.80
1:C:118:ARG:HH12	2:C:301:DTP:HN61	1.26	0.80
1:B:40:ARG:N	1:B:187:HIS:HD2	1.80	0.80
1:A:158:ASN:HD22	1:A:158:ASN:C	1.86	0.79
1:F:158:ASN:HD22	1:F:158:ASN:C	1.86	0.79
1:D:158:ASN:C	1:D:158:ASN:HD22	1.86	0.79
1:H:128:GLU:HA	1:H:131:LEU:HD12	1.64	0.79
1:C:40:ARG:N	1:C:187:HIS:HD2	1.79	0.79
1:C:158:ASN:C	1:C:158:ASN:HD22	1.86	0.79
1:D:128:GLU:HA	1:D:131:LEU:HD12	1.64	0.79
1:H:118:ARG:HH12	2:H:301:DTP:HN61	1.31	0.78
1:B:128:GLU:HA	1:B:131:LEU:HD12	1.64	0.78
1:G:40:ARG:N	1:G:187:HIS:HD2	1.80	0.78
1:C:127:PRO:O	1:C:131:LEU:HG	1.84	0.78
1:C:58:LEU:HA	1:C:267:MET:HE1	1.65	0.78
1:B:158:ASN:C	1:B:158:ASN:HD22	1.86	0.78
1:A:40:ARG:N	1:A:187:HIS:HD2	1.80	0.77
1:H:127:PRO:O	1:H:131:LEU:HG	1.84	0.77
1:B:58:LEU:HD23	1:B:267:MET:HE1	1.67	0.77
1:C:128:GLU:HA	1:C:131:LEU:HD12	1.64	0.77
1:F:127:PRO:O	1:F:131:LEU:HG	1.84	0.77
1:H:40:ARG:N	1:H:187:HIS:HD2	1.80	0.77
1:E:40:ARG:N	1:E:187:HIS:HD2	1.80	0.77
1:E:128:GLU:HA	1:E:131:LEU:HD12	1.65	0.77
1:H:158:ASN:HD22	1:H:158:ASN:C	1.86	0.77
1:A:127:PRO:O	1:A:131:LEU:HG	1.84	0.77
1:D:127:PRO:O	1:D:131:LEU:HG	1.84	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:40:ARG:N	1:D:187:HIS:HD2	1.80	0.76
1:F:128:GLU:HA	1:F:131:LEU:HD12	1.64	0.76
1:A:128:GLU:HA	1:A:131:LEU:HD12	1.64	0.76
1:E:158:ASN:HD22	1:E:158:ASN:C	1.86	0.76
1:G:128:GLU:HA	1:G:131:LEU:HD12	1.64	0.76
1:G:158:ASN:C	1:G:158:ASN:HD22	1.86	0.76
1:B:127:PRO:O	1:B:131:LEU:HG	1.84	0.76
1:F:40:ARG:N	1:F:187:HIS:HD2	1.79	0.76
1:G:127:PRO:O	1:G:131:LEU:HG	1.84	0.76
1:E:58:LEU:HD23	1:E:267:MET:HE1	1.67	0.76
1:C:78:ILE:HG13	1:D:164:ILE:HD12	1.67	0.75
1:E:127:PRO:O	1:E:131:LEU:HG	1.84	0.75
1:B:78:ILE:HD11	1:B:95:LEU:HB2	1.70	0.73
1:A:78:ILE:HD11	1:A:95:LEU:HB2	1.71	0.73
1:A:58:LEU:HD23	1:A:267:MET:HE3	1.70	0.73
1:D:78:ILE:HD11	1:D:95:LEU:HB2	1.71	0.73
1:C:78:ILE:HD11	1:C:95:LEU:HB2	1.70	0.73
1:E:78:ILE:HD11	1:E:95:LEU:HB2	1.71	0.73
1:F:78:ILE:HD11	1:F:95:LEU:HB2	1.71	0.72
1:H:78:ILE:HD11	1:H:95:LEU:HB2	1.70	0.72
1:G:78:ILE:HD11	1:G:95:LEU:HB2	1.70	0.71
1:B:40:ARG:HH11	1:B:40:ARG:HG3	1.55	0.71
1:F:40:ARG:HG3	1:F:40:ARG:HH11	1.56	0.71
1:G:128:GLU:HA	1:G:131:LEU:CD1	2.21	0.70
1:D:58:LEU:HD23	1:D:267:MET:HE3	1.73	0.70
1:C:40:ARG:HG3	1:C:40:ARG:HH11	1.55	0.70
1:E:128:GLU:HA	1:E:131:LEU:CD1	2.21	0.70
1:H:178:TRP:HB2	3:H:808:HOH:O	1.92	0.70
1:E:40:ARG:HH11	1:E:40:ARG:HG3	1.56	0.70
1:D:40:ARG:HH11	1:D:40:ARG:HG3	1.55	0.70
1:A:40:ARG:HG3	1:A:40:ARG:HH11	1.55	0.70
1:A:128:GLU:HA	1:A:131:LEU:CD1	2.21	0.70
1:H:40:ARG:HG3	1:H:40:ARG:HH11	1.55	0.70
1:H:128:GLU:HA	1:H:131:LEU:CD1	2.21	0.70
1:F:128:GLU:HA	1:F:131:LEU:CD1	2.21	0.70
1:D:128:GLU:HA	1:D:131:LEU:CD1	2.21	0.69
1:C:128:GLU:HA	1:C:131:LEU:CD1	2.21	0.69
1:D:58:LEU:HA	1:D:267:MET:HE1	1.74	0.69
1:G:178:TRP:HB2	3:G:708:HOH:O	1.92	0.69
1:G:58:LEU:HA	1:G:267:MET:HE1	1.74	0.69
1:B:128:GLU:HA	1:B:131:LEU:CD1	2.21	0.69
1:G:58:LEU:HD23	1:G:267:MET:HE3	1.74	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:178:TRP:HB2	3:F:608:HOH:O	1.92	0.69
1:G:40:ARG:HH11	1:G:40:ARG:HG3	1.56	0.69
1:A:178:TRP:HB2	3:A:307:HOH:O	1.92	0.69
1:B:178:TRP:HB2	3:B:309:HOH:O	1.92	0.68
1:D:178:TRP:HB2	3:D:309:HOH:O	1.92	0.68
1:E:178:TRP:HB2	3:E:508:HOH:O	1.92	0.68
1:C:178:TRP:HB2	3:C:308:HOH:O	1.92	0.67
1:H:57:LEU:HD22	1:H:267:MET:SD	2.35	0.67
1:E:57:LEU:HD22	1:E:267:MET:SD	2.35	0.67
1:B:57:LEU:HD22	1:B:267:MET:SD	2.35	0.67
1:F:57:LEU:HD22	1:F:267:MET:SD	2.35	0.67
1:A:70:GLU:HB2	1:A:141:GLU:HB2	1.77	0.67
1:G:57:LEU:HD22	1:G:267:MET:SD	2.35	0.67
1:E:99:MET:HG3	1:E:100:TYR:N	2.10	0.66
1:A:58:LEU:HA	1:A:267:MET:HE1	1.76	0.66
1:G:99:MET:HG3	1:G:100:TYR:N	2.10	0.66
1:C:70:GLU:HB2	1:C:141:GLU:HB2	1.77	0.66
1:A:57:LEU:HD22	1:A:267:MET:SD	2.35	0.66
1:D:57:LEU:HD22	1:D:267:MET:SD	2.35	0.66
1:D:70:GLU:HB2	1:D:141:GLU:HB2	1.77	0.66
1:F:70:GLU:HB2	1:F:141:GLU:HB2	1.77	0.66
1:G:70:GLU:HB2	1:G:141:GLU:HB2	1.77	0.66
1:C:57:LEU:HD22	1:C:267:MET:SD	2.35	0.66
1:E:70:GLU:HB2	1:E:141:GLU:HB2	1.77	0.66
1:B:99:MET:HG3	1:B:100:TYR:N	2.10	0.66
1:F:99:MET:HG3	1:F:100:TYR:N	2.10	0.66
1:H:99:MET:HG3	1:H:100:TYR:N	2.10	0.65
1:D:99:MET:HG3	1:D:100:TYR:N	2.10	0.65
1:H:70:GLU:HB2	1:H:141:GLU:HB2	1.77	0.65
1:E:58:LEU:HA	1:E:267:MET:HE2	1.79	0.65
1:A:99:MET:HG3	1:A:100:TYR:N	2.10	0.65
1:C:99:MET:HG3	1:C:100:TYR:N	2.10	0.64
1:B:70:GLU:HB2	1:B:141:GLU:HB2	1.77	0.64
1:B:58:LEU:HA	1:B:267:MET:HE2	1.79	0.63
1:D:158:ASN:ND2	1:D:158:ASN:C	2.53	0.62
1:E:158:ASN:ND2	1:E:158:ASN:C	2.53	0.62
1:F:158:ASN:ND2	1:F:158:ASN:C	2.53	0.62
1:B:158:ASN:C	1:B:158:ASN:ND2	2.53	0.62
1:H:158:ASN:ND2	1:H:158:ASN:C	2.53	0.61
1:G:158:ASN:C	1:G:158:ASN:ND2	2.53	0.61
1:C:158:ASN:ND2	1:C:158:ASN:C	2.53	0.61
1:H:62:TYR:OH	1:H:271:ASN:HB2	2.01	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:62:TYR:OH	1:C:271:ASN:HB2	2.01	0.61
1:A:158:ASN:ND2	1:A:158:ASN:C	2.53	0.61
1:A:62:TYR:OH	1:A:271:ASN:HB2	2.01	0.61
1:F:58:LEU:HA	1:F:267:MET:HE2	1.83	0.60
1:D:164:ILE:HG12	3:D:308:HOH:O	2.01	0.60
1:E:62:TYR:OH	1:E:271:ASN:HB2	2.01	0.60
1:H:164:ILE:HG12	3:H:807:HOH:O	2.01	0.60
1:G:62:TYR:OH	1:G:271:ASN:HB2	2.01	0.60
1:E:164:ILE:HG12	3:E:507:HOH:O	2.01	0.60
1:B:62:TYR:OH	1:B:271:ASN:HB2	2.01	0.60
1:G:164:ILE:HG12	3:G:707:HOH:O	2.01	0.60
1:F:62:TYR:OH	1:F:271:ASN:HB2	2.01	0.60
1:H:58:LEU:HA	1:H:267:MET:HE2	1.83	0.60
1:A:65:TRP:CH2	1:A:274:VAL:HG21	2.37	0.59
1:F:65:TRP:CH2	1:F:274:VAL:HG21	2.37	0.59
1:D:62:TYR:OH	1:D:271:ASN:HB2	2.01	0.59
1:D:65:TRP:CH2	1:D:274:VAL:HG21	2.37	0.59
1:A:164:ILE:HG12	3:A:306:HOH:O	2.01	0.59
1:B:164:ILE:HG12	3:B:308:HOH:O	2.01	0.59
1:C:195:SER:OG	1:C:198:VAL:HG23	2.03	0.59
1:C:65:TRP:CH2	1:C:274:VAL:HG21	2.37	0.59
1:E:65:TRP:CH2	1:E:274:VAL:HG21	2.37	0.59
1:F:164:ILE:HG12	3:F:607:HOH:O	2.01	0.59
1:G:70:GLU:OE1	1:G:71:PRO:HD2	2.03	0.59
1:C:164:ILE:HG12	3:C:307:HOH:O	2.01	0.59
1:G:65:TRP:CH2	1:G:274:VAL:HG21	2.37	0.59
1:H:65:TRP:CH2	1:H:274:VAL:HG21	2.37	0.58
1:E:195:SER:OG	1:E:198:VAL:HG23	2.03	0.58
1:B:78:ILE:HG12	1:B:93:GLY:O	2.03	0.58
1:A:70:GLU:OE1	1:A:71:PRO:HD2	2.03	0.58
1:F:70:GLU:OE1	1:F:71:PRO:HD2	2.03	0.58
1:F:78:ILE:HG12	1:F:93:GLY:O	2.03	0.58
1:H:70:GLU:OE1	1:H:71:PRO:HD2	2.03	0.58
1:B:65:TRP:CH2	1:B:274:VAL:HG21	2.37	0.58
1:F:195:SER:OG	1:F:198:VAL:HG23	2.03	0.58
1:C:78:ILE:HG12	1:C:93:GLY:O	2.03	0.58
1:C:70:GLU:OE1	1:C:71:PRO:HD2	2.03	0.58
1:A:78:ILE:HG12	1:A:93:GLY:O	2.03	0.58
1:B:195:SER:OG	1:B:198:VAL:HG23	2.03	0.58
1:E:70:GLU:OE1	1:E:71:PRO:HD2	2.03	0.58
1:D:78:ILE:HG12	1:D:93:GLY:O	2.03	0.58
1:H:78:ILE:HG12	1:H:93:GLY:O	2.03	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:195:SER:OG	1:D:198:VAL:HG23	2.03	0.58
1:D:40:ARG:NH1	1:D:40:ARG:HG3	2.19	0.58
1:B:70:GLU:OE1	1:B:71:PRO:HD2	2.03	0.58
1:G:40:ARG:HG3	1:G:40:ARG:NH1	2.19	0.57
1:E:40:ARG:NH1	1:E:40:ARG:HG3	2.19	0.57
1:G:78:ILE:HG12	1:G:93:GLY:O	2.03	0.57
1:A:195:SER:OG	1:A:198:VAL:HG23	2.03	0.57
1:A:40:ARG:NH1	1:A:40:ARG:HG3	2.19	0.57
1:D:70:GLU:OE1	1:D:71:PRO:HD2	2.03	0.57
1:H:195:SER:OG	1:H:198:VAL:HG23	2.03	0.57
1:C:40:ARG:HG3	1:C:40:ARG:NH1	2.19	0.57
1:G:78:ILE:HG13	1:H:164:ILE:HD12	1.86	0.57
1:G:195:SER:OG	1:G:198:VAL:HG23	2.03	0.57
1:E:78:ILE:HG12	1:E:93:GLY:O	2.03	0.57
1:E:166:TRP:O	1:E:170:GLN:HG3	2.05	0.57
1:H:40:ARG:HG3	1:H:40:ARG:NH1	2.19	0.56
1:B:166:TRP:O	1:B:170:GLN:HG3	2.05	0.56
1:C:166:TRP:O	1:C:170:GLN:HG3	2.05	0.56
1:G:166:TRP:O	1:G:170:GLN:HG3	2.05	0.56
1:D:145:TYR:CD2	1:D:242:LEU:HD22	2.41	0.56
1:G:145:TYR:CD2	1:G:242:LEU:HD22	2.41	0.56
1:B:61:THR:HB	1:B:267:MET:HE3	1.88	0.56
1:A:166:TRP:O	1:A:170:GLN:HG3	2.05	0.56
1:H:166:TRP:O	1:H:170:GLN:HG3	2.05	0.56
1:A:145:TYR:CD2	1:A:242:LEU:HD22	2.41	0.56
1:H:61:THR:HB	1:H:267:MET:HE3	1.88	0.56
1:A:62:TYR:O	1:A:65:TRP:HB2	2.06	0.56
1:G:62:TYR:O	1:G:65:TRP:HB2	2.06	0.56
1:D:62:TYR:O	1:D:65:TRP:HB2	2.06	0.55
1:B:145:TYR:CD2	1:B:242:LEU:HD22	2.41	0.55
1:F:145:TYR:CD2	1:F:242:LEU:HD22	2.41	0.55
1:B:153:LYS:O	1:B:157:GLU:HG3	2.07	0.55
1:D:166:TRP:O	1:D:170:GLN:HG3	2.05	0.55
1:F:166:TRP:O	1:F:170:GLN:HG3	2.05	0.55
1:E:62:TYR:O	1:E:65:TRP:HB2	2.07	0.55
1:B:62:TYR:O	1:B:65:TRP:HB2	2.06	0.55
1:E:153:LYS:O	1:E:157:GLU:HG3	2.07	0.55
1:E:61:THR:HB	1:E:267:MET:HE3	1.88	0.55
1:H:153:LYS:O	1:H:157:GLU:HG3	2.07	0.55
1:C:145:TYR:CD2	1:C:242:LEU:HD22	2.41	0.55
1:C:153:LYS:O	1:C:157:GLU:HG3	2.07	0.55
1:E:145:TYR:CD2	1:E:242:LEU:HD22	2.41	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:145:TYR:CD2	1:H:242:LEU:HD22	2.41	0.55
1:G:212:LYS:HG3	1:G:212:LYS:O	2.07	0.55
1:B:40:ARG:HG3	1:B:40:ARG:NH1	2.19	0.55
1:D:153:LYS:O	1:D:157:GLU:HG3	2.07	0.55
1:G:58:LEU:HA	1:G:267:MET:CE	2.37	0.55
1:B:58:LEU:HA	1:B:267:MET:CE	2.37	0.54
1:H:62:TYR:O	1:H:65:TRP:HB2	2.06	0.54
1:C:62:TYR:O	1:C:65:TRP:HB2	2.06	0.54
1:A:212:LYS:O	1:A:212:LYS:HG3	2.07	0.54
1:H:212:LYS:HG3	1:H:212:LYS:O	2.07	0.54
1:F:153:LYS:O	1:F:157:GLU:HG3	2.07	0.54
1:F:61:THR:HB	1:F:267:MET:HE3	1.88	0.54
1:F:40:ARG:NH1	1:F:40:ARG:HG3	2.19	0.54
1:B:61:THR:CB	1:B:267:MET:HE3	2.38	0.54
1:H:58:LEU:HA	1:H:267:MET:CE	2.37	0.54
1:E:58:LEU:HA	1:E:267:MET:CE	2.37	0.54
1:E:203:LEU:HD23	1:E:203:LEU:O	2.08	0.54
1:H:61:THR:CB	1:H:267:MET:HE3	2.38	0.54
1:C:212:LYS:HG3	1:C:212:LYS:O	2.07	0.54
1:B:212:LYS:HG3	1:B:212:LYS:O	2.07	0.54
1:E:212:LYS:O	1:E:212:LYS:HG3	2.07	0.54
1:F:212:LYS:HG3	1:F:212:LYS:O	2.07	0.54
1:A:61:THR:CB	1:A:267:MET:HE2	2.38	0.54
1:B:71:PRO:HG2	1:B:117:SER:OG	2.08	0.54
1:G:183:ARG:NH2	1:H:179:GLU:OE1	2.33	0.54
1:F:203:LEU:HD23	1:F:203:LEU:O	2.08	0.54
1:F:58:LEU:HA	1:F:267:MET:CE	2.37	0.54
1:F:62:TYR:O	1:F:65:TRP:HB2	2.06	0.54
1:D:212:LYS:HG3	1:D:212:LYS:O	2.07	0.54
1:A:58:LEU:HA	1:A:267:MET:CE	2.37	0.53
1:G:61:THR:CB	1:G:267:MET:HE2	2.38	0.53
1:G:67:VAL:HG22	1:G:139:ILE:HB	1.90	0.53
1:A:153:LYS:O	1:A:157:GLU:HG3	2.07	0.53
1:E:61:THR:CB	1:E:267:MET:HE3	2.38	0.53
1:C:67:VAL:HG22	1:C:139:ILE:HB	1.90	0.53
1:G:203:LEU:O	1:G:203:LEU:HD23	2.08	0.53
1:G:153:LYS:O	1:G:157:GLU:HG3	2.07	0.53
1:G:154:ASN:HD21	1:G:221:GLN:HE21	1.56	0.53
1:A:61:THR:HB	1:A:267:MET:HE2	1.89	0.53
1:D:61:THR:CB	1:D:267:MET:HE2	2.38	0.53
1:E:71:PRO:HG2	1:E:117:SER:OG	2.08	0.53
1:C:148:ARG:HD2	1:C:173:HIS:ND1	2.24	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:154:ASN:HD21	1:F:221:GLN:HE21	1.56	0.53
1:D:58:LEU:HA	1:D:267:MET:CE	2.37	0.53
1:D:61:THR:HB	1:D:267:MET:HE2	1.90	0.53
1:E:67:VAL:HG22	1:E:139:ILE:HB	1.90	0.53
1:A:154:ASN:HD21	1:A:221:GLN:HE21	1.56	0.53
1:D:148:ARG:HD2	1:D:173:HIS:ND1	2.24	0.53
1:A:71:PRO:HG2	1:A:117:SER:OG	2.08	0.53
1:D:71:PRO:HG2	1:D:117:SER:OG	2.08	0.53
1:F:67:VAL:HG22	1:F:139:ILE:HB	1.90	0.53
1:H:203:LEU:HD23	1:H:203:LEU:O	2.08	0.53
1:A:203:LEU:HD23	1:A:203:LEU:O	2.08	0.53
1:G:148:ARG:HD2	1:G:173:HIS:ND1	2.24	0.53
1:B:203:LEU:O	1:B:203:LEU:HD23	2.08	0.53
1:A:148:ARG:HD2	1:A:173:HIS:ND1	2.24	0.53
1:D:203:LEU:O	1:D:203:LEU:HD23	2.08	0.53
1:F:61:THR:CB	1:F:267:MET:HE3	2.38	0.53
1:H:71:PRO:HG2	1:H:117:SER:OG	2.08	0.53
1:B:148:ARG:HD2	1:B:173:HIS:ND1	2.24	0.53
1:E:148:ARG:HD2	1:E:173:HIS:ND1	2.24	0.53
1:H:67:VAL:HG22	1:H:139:ILE:HB	1.90	0.53
1:F:148:ARG:HD2	1:F:173:HIS:ND1	2.24	0.53
1:G:71:PRO:HG2	1:G:117:SER:OG	2.08	0.53
1:D:67:VAL:HG22	1:D:139:ILE:HB	1.90	0.52
1:C:154:ASN:HD21	1:C:221:GLN:HE21	1.56	0.52
1:F:39:ARG:HA	1:F:187:HIS:CD2	2.45	0.52
1:F:71:PRO:HG2	1:F:117:SER:OG	2.08	0.52
1:H:148:ARG:HD2	1:H:173:HIS:ND1	2.24	0.52
1:C:203:LEU:HD23	1:C:203:LEU:O	2.08	0.52
1:B:67:VAL:HG22	1:B:139:ILE:HB	1.90	0.52
1:H:39:ARG:HA	1:H:187:HIS:CD2	2.45	0.52
1:B:39:ARG:HA	1:B:187:HIS:CD2	2.45	0.52
1:E:154:ASN:HD21	1:E:221:GLN:HE21	1.56	0.52
1:G:177:LEU:HD22	1:G:184:ILE:HD13	1.92	0.52
1:D:154:ASN:HD21	1:D:221:GLN:HE21	1.56	0.52
1:H:217:ALA:O	1:H:221:GLN:HG3	2.10	0.52
1:G:61:THR:HB	1:G:267:MET:HE2	1.90	0.52
1:D:217:ALA:O	1:D:221:GLN:HG3	2.10	0.52
1:D:177:LEU:HD22	1:D:184:ILE:HD13	1.92	0.52
1:C:128:GLU:O	1:C:131:LEU:HB2	2.10	0.52
1:G:128:GLU:O	1:G:131:LEU:HB2	2.10	0.52
1:C:71:PRO:HG2	1:C:117:SER:OG	2.08	0.52
1:F:217:ALA:O	1:F:221:GLN:HG3	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:217:ALA:O	1:G:221:GLN:HG3	2.10	0.52
1:C:39:ARG:HA	1:C:187:HIS:CD2	2.45	0.52
1:D:39:ARG:HA	1:D:187:HIS:CD2	2.45	0.52
1:H:128:GLU:O	1:H:131:LEU:HB2	2.10	0.52
1:H:154:ASN:HD21	1:H:221:GLN:HE21	1.56	0.52
1:B:154:ASN:HD21	1:B:221:GLN:HE21	1.56	0.52
1:C:217:ALA:O	1:C:221:GLN:HG3	2.10	0.51
1:G:39:ARG:HA	1:G:187:HIS:CD2	2.45	0.51
1:A:67:VAL:HG22	1:A:139:ILE:HB	1.90	0.51
1:E:203:LEU:HD23	1:E:203:LEU:C	2.31	0.51
1:A:203:LEU:HD23	1:A:203:LEU:C	2.31	0.51
1:D:203:LEU:C	1:D:203:LEU:HD23	2.31	0.51
1:E:39:ARG:HA	1:E:187:HIS:CD2	2.45	0.51
1:A:39:ARG:HA	1:A:187:HIS:CD2	2.45	0.51
1:B:203:LEU:C	1:B:203:LEU:HD23	2.31	0.51
1:E:217:ALA:O	1:E:221:GLN:HG3	2.10	0.51
1:B:128:GLU:O	1:B:131:LEU:HB2	2.10	0.51
1:E:128:GLU:O	1:E:131:LEU:HB2	2.10	0.51
1:A:128:GLU:O	1:A:131:LEU:HB2	2.10	0.51
1:F:203:LEU:HD23	1:F:203:LEU:C	2.31	0.51
1:G:203:LEU:C	1:G:203:LEU:HD23	2.31	0.51
1:B:217:ALA:O	1:B:221:GLN:HG3	2.10	0.51
1:B:177:LEU:HD22	1:B:184:ILE:HD13	1.92	0.51
1:H:203:LEU:HD23	1:H:203:LEU:C	2.31	0.51
1:H:142:ARG:HD3	3:H:814:HOH:O	2.11	0.51
1:F:128:GLU:O	1:F:131:LEU:HB2	2.10	0.51
1:B:72:VAL:HA	1:B:75:TRP:CE3	2.46	0.51
1:H:72:VAL:HA	1:H:75:TRP:CE3	2.46	0.51
1:C:203:LEU:C	1:C:203:LEU:HD23	2.31	0.51
1:E:177:LEU:HD22	1:E:184:ILE:HD13	1.92	0.51
1:G:72:VAL:HA	1:G:75:TRP:CE3	2.46	0.50
1:D:128:GLU:O	1:D:131:LEU:HB2	2.10	0.50
1:H:177:LEU:HD22	1:H:184:ILE:HD13	1.92	0.50
1:C:177:LEU:HD22	1:C:184:ILE:HD13	1.92	0.50
1:A:177:LEU:HD22	1:A:184:ILE:HD13	1.92	0.50
1:E:72:VAL:HA	1:E:75:TRP:CE3	2.46	0.50
1:C:72:VAL:HA	1:C:75:TRP:CE3	2.46	0.50
1:E:259:GLU:N	1:E:259:GLU:OE1	2.45	0.50
1:H:259:GLU:OE1	1:H:259:GLU:N	2.45	0.50
1:A:217:ALA:O	1:A:221:GLN:HG3	2.10	0.50
1:H:49:VAL:HG23	1:H:194:ALA:HB3	1.93	0.50
1:A:72:VAL:HA	1:A:75:TRP:CE3	2.46	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:58:LEU:HA	1:C:267:MET:CE	2.37	0.50
1:G:49:VAL:HG23	1:G:194:ALA:HB3	1.93	0.50
1:B:211:GLU:O	1:B:213:GLY:N	2.45	0.50
1:G:259:GLU:OE1	1:G:259:GLU:N	2.45	0.50
1:B:259:GLU:OE1	1:B:259:GLU:N	2.45	0.50
1:F:72:VAL:HA	1:F:75:TRP:CE3	2.46	0.50
1:G:58:LEU:HD23	1:G:267:MET:CE	2.41	0.50
1:A:211:GLU:O	1:A:213:GLY:N	2.45	0.50
1:E:211:GLU:O	1:E:213:GLY:N	2.45	0.50
1:F:177:LEU:HD22	1:F:184:ILE:HD13	1.92	0.49
1:A:142:ARG:HD3	3:A:313:HOH:O	2.11	0.49
1:D:72:VAL:HA	1:D:75:TRP:CE3	2.46	0.49
1:A:58:LEU:HD23	1:A:267:MET:CE	2.41	0.49
1:D:211:GLU:O	1:D:213:GLY:N	2.45	0.49
1:C:211:GLU:O	1:C:213:GLY:N	2.45	0.49
1:F:142:ARG:HD3	3:F:614:HOH:O	2.11	0.49
1:C:49:VAL:HG23	1:C:194:ALA:HB3	1.93	0.49
1:G:142:ARG:HD3	3:G:714:HOH:O	2.11	0.49
1:G:211:GLU:O	1:G:213:GLY:N	2.45	0.49
1:E:58:LEU:HD23	1:E:267:MET:CE	2.41	0.49
1:A:259:GLU:OE1	1:A:259:GLU:N	2.45	0.49
1:D:58:LEU:HD23	1:D:267:MET:CE	2.41	0.49
1:B:49:VAL:HG23	1:B:194:ALA:HB3	1.93	0.49
1:C:142:ARG:HD3	3:C:314:HOH:O	2.11	0.49
1:F:211:GLU:O	1:F:213:GLY:N	2.45	0.49
1:C:259:GLU:N	1:C:259:GLU:OE1	2.45	0.49
1:H:211:GLU:O	1:H:213:GLY:N	2.45	0.49
1:D:142:ARG:HD3	3:D:315:HOH:O	2.11	0.49
1:B:142:ARG:HD3	3:B:315:HOH:O	2.11	0.49
1:F:49:VAL:HG23	1:F:194:ALA:HB3	1.93	0.49
1:F:259:GLU:N	1:F:259:GLU:OE1	2.45	0.49
1:D:259:GLU:OE1	1:D:259:GLU:N	2.45	0.49
1:E:49:VAL:HG23	1:E:194:ALA:HB3	1.93	0.49
1:E:164:ILE:HD12	1:F:78:ILE:HG13	1.95	0.49
1:D:49:VAL:HG23	1:D:194:ALA:HB3	1.93	0.49
1:B:58:LEU:HD23	1:B:267:MET:CE	2.41	0.48
1:G:164:ILE:HD12	1:H:78:ILE:HG13	1.94	0.48
1:A:49:VAL:HG23	1:A:194:ALA:HB3	1.93	0.48
1:E:142:ARG:HD3	3:E:514:HOH:O	2.11	0.48
1:F:200:LEU:HD22	1:F:216:LEU:HB2	1.96	0.48
1:A:200:LEU:HD11	1:A:204:TYR:CE1	2.49	0.48
1:H:45:GLY:O	1:H:142:ARG:NH1	2.47	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:200:LEU:HD22	1:D:216:LEU:HB2	1.96	0.48
1:D:200:LEU:HD11	1:D:204:TYR:CE1	2.49	0.48
1:B:200:LEU:HD11	1:B:204:TYR:CE1	2.49	0.48
1:D:45:GLY:O	1:D:142:ARG:NH1	2.47	0.48
1:B:200:LEU:HD22	1:B:216:LEU:HB2	1.96	0.48
1:E:200:LEU:HD11	1:E:204:TYR:CE1	2.49	0.48
1:G:200:LEU:HD22	1:G:216:LEU:HB2	1.96	0.48
1:F:200:LEU:HD11	1:F:204:TYR:CE1	2.49	0.48
1:G:45:GLY:O	1:G:142:ARG:NH1	2.47	0.48
1:C:45:GLY:O	1:C:142:ARG:NH1	2.47	0.48
1:E:45:GLY:O	1:E:142:ARG:NH1	2.47	0.47
1:A:45:GLY:O	1:A:142:ARG:NH1	2.47	0.47
1:H:200:LEU:HD11	1:H:204:TYR:CE1	2.49	0.47
1:C:200:LEU:HD11	1:C:204:TYR:CE1	2.49	0.47
1:E:200:LEU:HD22	1:E:216:LEU:HB2	1.96	0.47
1:C:200:LEU:HD22	1:C:216:LEU:HB2	1.96	0.47
1:B:45:GLY:O	1:B:142:ARG:NH1	2.47	0.47
1:F:233:LYS:HB2	1:G:232:HIS:CG	2.50	0.47
1:E:65:TRP:HH2	1:E:274:VAL:HG21	1.80	0.47
1:A:209:GLU:O	1:A:212:LYS:HG2	2.15	0.47
1:E:209:GLU:O	1:E:212:LYS:HG2	2.15	0.47
1:A:200:LEU:HD22	1:A:216:LEU:HB2	1.96	0.47
1:G:200:LEU:HD11	1:G:204:TYR:CE1	2.49	0.47
1:H:200:LEU:HD22	1:H:216:LEU:HB2	1.96	0.47
1:D:240:GLU:O	1:D:243:MET:HB2	2.15	0.47
1:F:45:GLY:O	1:F:142:ARG:NH1	2.47	0.47
1:F:92:LEU:N	3:F:624:HOH:O	2.48	0.47
1:A:164:ILE:HD12	1:B:78:ILE:HG13	1.96	0.47
1:H:209:GLU:O	1:H:212:LYS:HG2	2.15	0.47
1:F:209:GLU:O	1:F:212:LYS:HG2	2.15	0.47
1:F:240:GLU:O	1:F:243:MET:HB2	2.15	0.47
1:D:209:GLU:O	1:D:212:LYS:HG2	2.15	0.47
1:G:41:LEU:HB2	1:G:139:ILE:HD13	1.97	0.47
1:B:240:GLU:O	1:B:243:MET:HB2	2.15	0.47
1:A:78:ILE:HG13	1:B:164:ILE:HD12	1.96	0.46
1:C:209:GLU:O	1:C:212:LYS:HG2	2.15	0.46
1:C:65:TRP:HH2	1:C:274:VAL:HG21	1.80	0.46
1:G:209:GLU:O	1:G:212:LYS:HG2	2.15	0.46
1:E:41:LEU:HB2	1:E:139:ILE:HD13	1.97	0.46
1:H:41:LEU:HB2	1:H:139:ILE:HD13	1.97	0.46
1:C:240:GLU:O	1:C:243:MET:HB2	2.15	0.46
1:A:158:ASN:ND2	1:A:160:SER:H	2.14	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:158:ASN:ND2	1:H:160:SER:H	2.14	0.46
1:G:65:TRP:HH2	1:G:274:VAL:HG21	1.80	0.46
1:A:240:GLU:O	1:A:243:MET:HB2	2.15	0.46
1:F:158:ASN:ND2	1:F:160:SER:H	2.14	0.46
1:A:70:GLU:HB2	1:A:141:GLU:CB	2.45	0.46
1:H:240:GLU:O	1:H:243:MET:HB2	2.15	0.46
1:G:240:GLU:O	1:G:243:MET:HB2	2.15	0.46
1:H:92:LEU:N	3:H:824:HOH:O	2.48	0.46
1:B:41:LEU:HB2	1:B:139:ILE:HD13	1.97	0.46
1:G:92:LEU:N	3:G:724:HOH:O	2.48	0.46
1:B:158:ASN:ND2	1:B:160:SER:H	2.14	0.46
1:B:209:GLU:O	1:B:212:LYS:HG2	2.15	0.46
1:C:92:LEU:N	3:C:324:HOH:O	2.48	0.46
1:D:92:LEU:N	3:D:325:HOH:O	2.48	0.46
1:E:240:GLU:O	1:E:243:MET:HB2	2.15	0.46
1:G:245:ILE:HG23	1:G:246:PRO:HD2	1.98	0.46
1:D:41:LEU:HB2	1:D:139:ILE:HD13	1.97	0.46
1:C:41:LEU:HB2	1:C:139:ILE:HD13	1.97	0.46
1:F:41:LEU:HB2	1:F:139:ILE:HD13	1.97	0.46
1:B:92:LEU:N	3:B:325:HOH:O	2.48	0.46
1:B:210:GLU:H	1:B:210:GLU:CD	2.20	0.46
1:A:92:LEU:N	3:A:323:HOH:O	2.48	0.46
1:E:245:ILE:HG23	1:E:246:PRO:HD2	1.98	0.45
1:A:245:ILE:HG23	1:A:246:PRO:HD2	1.98	0.45
1:D:158:ASN:ND2	1:D:160:SER:H	2.14	0.45
1:H:70:GLU:HB2	1:H:141:GLU:CB	2.45	0.45
1:D:200:LEU:HD11	1:D:204:TYR:HE1	1.81	0.45
1:G:200:LEU:HD11	1:G:204:TYR:HE1	1.81	0.45
1:F:210:GLU:CD	1:F:210:GLU:H	2.20	0.45
1:E:158:ASN:ND2	1:E:160:SER:H	2.14	0.45
1:C:158:ASN:ND2	1:C:160:SER:H	2.14	0.45
1:A:210:GLU:H	1:A:210:GLU:CD	2.20	0.45
1:C:210:GLU:CD	1:C:210:GLU:H	2.20	0.45
1:E:92:LEU:N	3:E:524:HOH:O	2.48	0.45
1:G:210:GLU:CD	1:G:210:GLU:H	2.20	0.45
1:E:210:GLU:CD	1:E:210:GLU:H	2.20	0.45
1:A:118:ARG:O	1:A:121:VAL:HG12	2.17	0.45
1:A:65:TRP:HH2	1:A:274:VAL:HG21	1.80	0.45
1:D:65:TRP:HH2	1:D:274:VAL:HG21	1.80	0.45
1:B:200:LEU:HD11	1:B:204:TYR:HE1	1.81	0.45
1:H:210:GLU:H	1:H:210:GLU:CD	2.20	0.45
1:F:245:ILE:HG23	1:F:246:PRO:HD2	1.98	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:118:ARG:O	1:H:121:VAL:HG12	2.17	0.45
1:C:70:GLU:HB2	1:C:141:GLU:CB	2.46	0.45
1:E:183:ARG:NH2	1:F:179:GLU:OE1	2.40	0.45
1:E:179:GLU:OE1	1:F:183:ARG:NH2	2.34	0.45
1:G:158:ASN:ND2	1:G:160:SER:H	2.14	0.45
1:A:41:LEU:HB2	1:A:139:ILE:HD13	1.97	0.45
1:E:118:ARG:O	1:E:121:VAL:HG12	2.17	0.45
1:B:245:ILE:HG23	1:B:246:PRO:HD2	1.98	0.45
1:G:118:ARG:O	1:G:121:VAL:HG12	2.17	0.45
1:D:245:ILE:HG23	1:D:246:PRO:HD2	1.98	0.45
1:E:200:LEU:HD11	1:E:204:TYR:HE1	1.81	0.45
1:H:200:LEU:HD11	1:H:204:TYR:HE1	1.82	0.45
1:C:245:ILE:HG23	1:C:246:PRO:HD2	1.98	0.44
1:F:70:GLU:HB2	1:F:141:GLU:CB	2.45	0.44
1:H:162:SER:OG	1:H:165:GLU:HG3	2.17	0.44
1:A:46:ASN:OD1	1:A:47:ILE:N	2.48	0.44
1:F:118:ARG:O	1:F:121:VAL:HG12	2.17	0.44
1:G:162:SER:OG	1:G:165:GLU:HG3	2.17	0.44
1:H:245:ILE:HG23	1:H:246:PRO:HD2	1.98	0.44
1:F:65:TRP:HH2	1:F:274:VAL:HG21	1.80	0.44
1:C:162:SER:OG	1:C:165:GLU:HG3	2.17	0.44
1:F:162:SER:OG	1:F:165:GLU:HG3	2.17	0.44
1:C:118:ARG:O	1:C:121:VAL:HG12	2.17	0.44
1:A:200:LEU:HD11	1:A:204:TYR:HE1	1.81	0.44
1:E:162:SER:OG	1:E:165:GLU:HG3	2.18	0.44
1:B:118:ARG:O	1:B:121:VAL:HG12	2.17	0.44
1:F:46:ASN:OD1	1:F:47:ILE:N	2.48	0.44
1:B:162:SER:OG	1:B:165:GLU:HG3	2.17	0.44
1:D:118:ARG:O	1:D:121:VAL:HG12	2.17	0.44
1:D:162:SER:OG	1:D:165:GLU:HG3	2.17	0.44
1:B:70:GLU:HA	1:B:71:PRO:HD3	1.76	0.44
1:A:162:SER:OG	1:A:165:GLU:HG3	2.17	0.44
1:D:210:GLU:H	1:D:210:GLU:CD	2.20	0.44
1:A:172:TRP:CZ2	1:B:116:LEU:HB2	2.53	0.44
1:F:72:VAL:HG21	2:F:301:DTP:O2A	2.17	0.44
1:C:46:ASN:OD1	1:C:47:ILE:N	2.48	0.44
1:A:70:GLU:HG2	3:A:321:HOH:O	2.18	0.43
1:D:70:GLU:HB2	1:D:141:GLU:CB	2.45	0.43
1:H:46:ASN:OD1	1:H:47:ILE:N	2.48	0.43
1:C:158:ASN:HD22	1:C:159:GLY:N	2.17	0.43
1:D:70:GLU:HG2	3:D:323:HOH:O	2.18	0.43
1:F:70:GLU:HG2	3:F:622:HOH:O	2.18	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:200:LEU:HD11	1:F:204:TYR:HE1	1.81	0.43
1:E:70:GLU:HB2	1:E:141:GLU:CB	2.45	0.43
1:G:179:GLU:OE1	1:H:183:ARG:NH2	2.39	0.43
1:D:256:PHE:CD1	1:D:256:PHE:C	2.92	0.43
1:H:58:LEU:HD23	1:H:267:MET:CE	2.41	0.43
1:D:72:VAL:HG21	2:D:301:DTP:O2A	2.19	0.43
1:H:256:PHE:C	1:H:256:PHE:CD1	2.92	0.43
1:C:70:GLU:HA	1:C:71:PRO:HD3	1.76	0.43
1:B:256:PHE:C	1:B:256:PHE:CD1	2.92	0.43
1:B:65:TRP:HH2	1:B:274:VAL:HG21	1.80	0.43
1:E:46:ASN:OD1	1:E:47:ILE:N	2.48	0.43
1:F:197:GLN:H	1:F:197:GLN:CD	2.22	0.43
1:C:256:PHE:C	1:C:256:PHE:CD1	2.92	0.43
1:E:158:ASN:HD22	1:E:159:GLY:N	2.17	0.42
1:G:70:GLU:HA	1:G:71:PRO:HD3	1.76	0.42
1:A:116:LEU:HB2	1:B:172:TRP:CZ2	2.54	0.42
1:A:197:GLN:H	1:A:197:GLN:CD	2.22	0.42
1:G:70:GLU:HG2	3:G:722:HOH:O	2.18	0.42
1:H:70:GLU:HG2	3:H:822:HOH:O	2.18	0.42
1:H:65:TRP:HH2	1:H:274:VAL:HG21	1.80	0.42
1:C:108:TYR:O	1:C:112:THR:HG23	2.20	0.42
2:H:301:DTP:H5'2	2:H:301:DTP:H2'2	1.81	0.42
1:B:158:ASN:HD22	1:B:159:GLY:N	2.17	0.42
1:G:158:ASN:HD22	1:G:159:GLY:N	2.17	0.42
1:B:70:GLU:HG2	3:B:323:HOH:O	2.18	0.42
1:D:135:LYS:HE3	1:D:275:LYS:HE2	2.02	0.42
1:G:108:TYR:O	1:G:112:THR:HG23	2.20	0.42
1:D:276:ASN:HD22	1:D:276:ASN:N	2.17	0.42
1:F:58:LEU:HD23	1:F:267:MET:CE	2.41	0.42
1:F:158:ASN:HD22	1:F:159:GLY:N	2.17	0.42
1:C:70:GLU:HG2	3:C:322:HOH:O	2.18	0.42
1:E:256:PHE:C	1:E:256:PHE:CD1	2.92	0.42
1:A:256:PHE:C	1:A:256:PHE:CD1	2.92	0.42
1:F:276:ASN:HD22	1:F:276:ASN:N	2.17	0.42
1:B:70:GLU:HB2	1:B:141:GLU:CB	2.45	0.42
1:C:200:LEU:HD11	1:C:204:TYR:HE1	1.82	0.42
1:B:197:GLN:H	1:B:197:GLN:CD	2.22	0.42
1:G:135:LYS:HE3	1:G:275:LYS:HE2	2.02	0.42
1:D:128:GLU:CD	1:D:128:GLU:H	2.23	0.42
1:E:108:TYR:O	1:E:112:THR:HG23	2.20	0.42
1:H:197:GLN:H	1:H:197:GLN:CD	2.22	0.42
1:G:72:VAL:HG21	2:G:301:DTP:O2A	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:158:ASN:HD22	1:A:159:GLY:N	2.17	0.42
1:E:70:GLU:HG2	3:E:522:HOH:O	2.18	0.42
1:E:135:LYS:HE3	1:E:275:LYS:HE2	2.02	0.42
1:F:256:PHE:CD1	1:F:256:PHE:C	2.92	0.42
1:H:108:TYR:O	1:H:112:THR:HG23	2.20	0.42
1:H:276:ASN:HD22	1:H:276:ASN:N	2.17	0.42
1:E:128:GLU:H	1:E:128:GLU:CD	2.23	0.42
1:G:128:GLU:CD	1:G:128:GLU:H	2.23	0.42
1:C:135:LYS:HE3	1:C:275:LYS:HE2	2.02	0.42
1:H:135:LYS:HE3	1:H:275:LYS:HE2	2.01	0.42
1:E:187:HIS:O	1:E:246:PRO:HD2	2.20	0.42
1:B:72:VAL:HG21	2:B:301:DTP:O2A	2.20	0.42
1:B:75:TRP:HD1	1:B:113:PHE:HE2	1.68	0.42
1:F:128:GLU:CD	1:F:128:GLU:H	2.23	0.42
1:A:75:TRP:HD1	1:A:113:PHE:HE2	1.68	0.42
1:A:108:TYR:O	1:A:112:THR:HG23	2.20	0.42
1:C:197:GLN:CD	1:C:197:GLN:H	2.22	0.42
1:G:256:PHE:CD1	1:G:256:PHE:C	2.92	0.42
1:C:276:ASN:HD22	1:C:276:ASN:N	2.17	0.42
1:D:158:ASN:HD22	1:D:159:GLY:N	2.17	0.42
1:F:108:TYR:O	1:F:112:THR:HG23	2.20	0.42
1:B:108:TYR:O	1:B:112:THR:HG23	2.20	0.42
1:H:187:HIS:O	1:H:246:PRO:HD2	2.20	0.41
1:B:187:HIS:O	1:B:246:PRO:HD2	2.20	0.41
2:G:301:DTP:H2'2	2:G:301:DTP:H5'2	1.83	0.41
1:D:75:TRP:HD1	1:D:113:PHE:HE2	1.68	0.41
1:H:128:GLU:CD	1:H:128:GLU:H	2.23	0.41
1:E:78:ILE:HG13	1:F:164:ILE:HD12	2.00	0.41
1:D:108:TYR:O	1:D:112:THR:HG23	2.20	0.41
1:G:46:ASN:OD1	1:G:47:ILE:N	2.48	0.41
1:D:197:GLN:H	1:D:197:GLN:CD	2.22	0.41
1:C:187:HIS:O	1:C:246:PRO:HD2	2.20	0.41
1:A:276:ASN:HD22	1:A:276:ASN:N	2.17	0.41
1:G:197:GLN:CD	1:G:197:GLN:H	2.22	0.41
1:H:158:ASN:HD22	1:H:159:GLY:N	2.17	0.41
1:F:70:GLU:HA	1:F:71:PRO:HD3	1.76	0.41
1:E:197:GLN:H	1:E:197:GLN:CD	2.22	0.41
1:G:276:ASN:N	1:G:276:ASN:HD22	2.17	0.41
1:A:135:LYS:HE3	1:A:275:LYS:HE2	2.02	0.41
1:C:128:GLU:H	1:C:128:GLU:CD	2.23	0.41
1:A:128:GLU:CD	1:A:128:GLU:H	2.23	0.41
1:F:177:LEU:HD22	1:F:184:ILE:CD1	2.51	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:135:LYS:HE3	1:F:275:LYS:HE2	2.02	0.41
1:E:276:ASN:N	1:E:276:ASN:HD22	2.17	0.41
1:B:124:GLU:HA	1:B:125:PRO:HD3	1.91	0.41
1:G:187:HIS:O	1:G:246:PRO:HD2	2.20	0.41
1:H:75:TRP:HD1	1:H:113:PHE:HE2	1.68	0.41
1:A:70:GLU:HA	1:A:71:PRO:HD3	1.76	0.41
1:F:187:HIS:O	1:F:246:PRO:HD2	2.20	0.41
1:B:244:ASN:ND2	1:E:134:ARG:NH2	2.68	0.41
1:G:75:TRP:HD1	1:G:113:PHE:HE2	1.68	0.41
2:B:301:DTP:H5'2	2:B:301:DTP:H2'2	1.80	0.41
1:B:46:ASN:OD1	1:B:47:ILE:N	2.48	0.41
1:C:75:TRP:HD1	1:C:113:PHE:HE2	1.68	0.41
1:H:177:LEU:HD22	1:H:184:ILE:CD1	2.51	0.41
1:D:187:HIS:O	1:D:246:PRO:HD2	2.20	0.41
1:B:128:GLU:CD	1:B:128:GLU:H	2.23	0.41
1:H:64:GLU:O	1:H:129:LYS:HE2	2.21	0.41
1:B:135:LYS:HE3	1:B:275:LYS:HE2	2.02	0.41
1:H:72:VAL:HG21	2:H:301:DTP:O2A	2.21	0.41
1:A:172:TRP:CE2	1:B:116:LEU:HB2	2.56	0.41
1:E:75:TRP:HD1	1:E:113:PHE:HE2	1.68	0.40
1:B:276:ASN:HD22	1:B:276:ASN:N	2.17	0.40
1:A:177:LEU:HD22	1:A:184:ILE:CD1	2.51	0.40
1:G:64:GLU:O	1:G:129:LYS:HE2	2.21	0.40
1:A:187:HIS:O	1:A:246:PRO:HD2	2.20	0.40
1:C:72:VAL:HG21	2:C:301:DTP:O2A	2.21	0.40
1:G:70:GLU:HB2	1:G:141:GLU:CB	2.45	0.40
1:E:70:GLU:HA	1:E:71:PRO:HD3	1.76	0.40
1:F:64:GLU:O	1:F:129:LYS:HE2	2.21	0.40
1:B:75:TRP:HD1	1:B:113:PHE:CE2	2.40	0.40
1:H:128:GLU:HA	1:H:131:LEU:CG	2.52	0.40
1:C:177:LEU:HD22	1:C:184:ILE:CD1	2.51	0.40
1:F:233:LYS:HG3	1:G:232:HIS:HB3	2.04	0.40
1:C:64:GLU:O	1:C:129:LYS:HE2	2.21	0.40
1:D:177:LEU:HD22	1:D:184:ILE:CD1	2.51	0.40
1:E:177:LEU:HD22	1:E:184:ILE:CD1	2.51	0.40
1:D:46:ASN:OD1	1:D:47:ILE:N	2.48	0.40

All (20) 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:C:135:LYS:CD	1:E:240:GLU:OE1[2_756]	0.72	1.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:135:LYS:CE	1:E:240:GLU:OE1[2.756]	0.83	1.37
1:C:135:LYS:CE	1:E:240:GLU:CD[2.756]	1.09	1.11
1:F:129:LYS:CB	1:G:63:PRO:CB[1.655]	1.21	0.99
1:B:131:LEU:CD2	1:H:275:LYS:NZ[1.565]	1.48	0.72
1:C:276:ASN:OD1	1:E:185:THR:CG2[2.756]	1.74	0.46
1:C:135:LYS:CD	1:E:240:GLU:CD[2.756]	1.81	0.39
1:F:129:LYS:CD	1:G:63:PRO:O[1.655]	1.85	0.35
1:D:178:TRP:NE1	1:G:134:ARG:NH1[2.656]	1.88	0.32
1:B:123:LEU:O	1:H:134:ARG:CD[1.565]	1.89	0.31
1:C:135:LYS:CE	1:E:240:GLU:OE2[2.756]	1.91	0.29
1:A:135:LYS:NZ	1:G:240:GLU:OE2[1.565]	1.97	0.23
1:C:135:LYS:CG	1:E:240:GLU:OE1[2.756]	1.98	0.22
1:C:135:LYS:NZ	1:E:240:GLU:OE1[2.756]	2.00	0.20
1:D:178:TRP:CD1	1:G:134:ARG:NH1[2.656]	2.05	0.15
1:F:129:LYS:CG	1:G:63:PRO:CB[1.655]	2.08	0.12
1:B:122:GLN:O	1:H:134:ARG:NH1[1.565]	2.15	0.05
1:E:101:ARG:NH2	1:G:92:LEU:CD1[2.656]	2.15	0.05
1:C:135:LYS:NZ	1:E:240:GLU:CD[2.756]	2.18	0.02
1:C:276:ASN:O	1:E:182:SER:CA[2.756]	2.19	0.01

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	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
1	B	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
1	C	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
1	D	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
1	E	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
1	F	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
1	G	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
1	H	225/241 (93%)	203 (90%)	18 (8%)	4 (2%)	13	39
All	All	1800/1928 (93%)	1624 (90%)	144 (8%)	32 (2%)	13	39

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	71	PRO
1	A	207	ALA
1	B	71	PRO
1	B	207	ALA
1	C	71	PRO
1	C	207	ALA
1	D	71	PRO
1	D	207	ALA
1	E	71	PRO
1	E	207	ALA
1	F	71	PRO
1	F	207	ALA
1	G	71	PRO
1	G	207	ALA
1	H	71	PRO
1	H	207	ALA
1	A	212	LYS
1	A	255	ASP
1	B	212	LYS
1	B	255	ASP
1	C	212	LYS
1	C	255	ASP
1	D	212	LYS
1	D	255	ASP
1	E	212	LYS
1	E	255	ASP
1	F	212	LYS
1	F	255	ASP
1	G	212	LYS
1	G	255	ASP
1	H	212	LYS
1	H	255	ASP

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	209/216 (97%)	196 (94%)	13 (6%)	26	60
1	B	209/216 (97%)	196 (94%)	13 (6%)	26	60
1	C	209/216 (97%)	196 (94%)	13 (6%)	26	60
1	D	209/216 (97%)	196 (94%)	13 (6%)	26	60
1	E	209/216 (97%)	196 (94%)	13 (6%)	26	60
1	F	209/216 (97%)	196 (94%)	13 (6%)	26	60
1	G	209/216 (97%)	196 (94%)	13 (6%)	26	60
1	H	209/216 (97%)	196 (94%)	13 (6%)	26	60
All	All	1672/1728 (97%)	1568 (94%)	104 (6%)	26	60

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

Mol	Chain	Res	Type
1	A	39	ARG
1	A	95	LEU
1	A	99	MET
1	A	111	GLN
1	A	122	GLN
1	A	144	VAL
1	A	158	ASN
1	A	171	ASP
1	A	197	GLN
1	A	208	ARG
1	A	258	GLU
1	A	259	GLU
1	A	266	LEU
1	B	39	ARG
1	B	95	LEU
1	B	99	MET
1	B	111	GLN
1	B	122	GLN
1	B	144	VAL
1	B	158	ASN
1	B	171	ASP
1	B	197	GLN
1	B	208	ARG
1	B	258	GLU
1	B	259	GLU
1	B	266	LEU
1	C	39	ARG

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Mol	Chain	Res	Type
1	C	95	LEU
1	C	99	MET
1	C	111	GLN
1	C	122	GLN
1	C	144	VAL
1	C	158	ASN
1	C	171	ASP
1	C	197	GLN
1	C	208	ARG
1	C	258	GLU
1	C	259	GLU
1	C	266	LEU
1	D	39	ARG
1	D	95	LEU
1	D	99	MET
1	D	111	GLN
1	D	122	GLN
1	D	144	VAL
1	D	158	ASN
1	D	171	ASP
1	D	197	GLN
1	D	208	ARG
1	D	258	GLU
1	D	259	GLU
1	D	266	LEU
1	E	39	ARG
1	E	95	LEU
1	E	99	MET
1	E	111	GLN
1	E	122	GLN
1	E	144	VAL
1	E	158	ASN
1	E	171	ASP
1	E	197	GLN
1	E	208	ARG
1	E	258	GLU
1	E	259	GLU
1	E	266	LEU
1	F	39	ARG
1	F	95	LEU
1	F	99	MET
1	F	111	GLN

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Mol	Chain	Res	Type
1	F	122	GLN
1	F	144	VAL
1	F	158	ASN
1	F	171	ASP
1	F	197	GLN
1	F	208	ARG
1	F	258	GLU
1	F	259	GLU
1	F	266	LEU
1	G	39	ARG
1	G	95	LEU
1	G	99	MET
1	G	111	GLN
1	G	122	GLN
1	G	144	VAL
1	G	158	ASN
1	G	171	ASP
1	G	197	GLN
1	G	208	ARG
1	G	258	GLU
1	G	259	GLU
1	G	266	LEU
1	H	39	ARG
1	H	95	LEU
1	H	99	MET
1	H	111	GLN
1	H	122	GLN
1	H	144	VAL
1	H	158	ASN
1	H	171	ASP
1	H	197	GLN
1	H	208	ARG
1	H	258	GLU
1	H	259	GLU
1	H	266	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	A	187	HIS
1	A	221	GLN

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Mol	Chain	Res	Type
1	A	271	ASN
1	A	276	ASN
1	B	158	ASN
1	B	187	HIS
1	B	221	GLN
1	B	271	ASN
1	B	276	ASN
1	C	158	ASN
1	C	187	HIS
1	C	221	GLN
1	C	271	ASN
1	C	276	ASN
1	D	158	ASN
1	D	187	HIS
1	D	221	GLN
1	D	271	ASN
1	D	276	ASN
1	E	158	ASN
1	E	187	HIS
1	E	221	GLN
1	E	271	ASN
1	E	276	ASN
1	F	158	ASN
1	F	187	HIS
1	F	221	GLN
1	F	271	ASN
1	F	276	ASN
1	G	158	ASN
1	G	187	HIS
1	G	221	GLN
1	G	271	ASN
1	G	276	ASN
1	H	158	ASN
1	H	187	HIS
1	H	221	GLN
1	H	271	ASN
1	H	276	ASN

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 ⓘ

8 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	DTP	A	301	-	32,32,32	0.91	0	50,50,50	1.93	8 (16%)
2	DTP	B	301	-	32,32,32	0.91	0	50,50,50	1.97	7 (14%)
2	DTP	C	301	-	32,32,32	0.88	1 (3%)	50,50,50	2.03	8 (16%)
2	DTP	D	301	-	32,32,32	0.98	2 (6%)	50,50,50	2.01	7 (14%)
2	DTP	E	301	-	32,32,32	0.97	1 (3%)	50,50,50	1.98	9 (18%)
2	DTP	F	301	-	32,32,32	0.91	1 (3%)	50,50,50	2.07	7 (14%)
2	DTP	G	301	-	32,32,32	0.92	1 (3%)	50,50,50	2.03	7 (14%)
2	DTP	H	301	-	32,32,32	0.90	1 (3%)	50,50,50	1.99	6 (12%)

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	DTP	A	301	-	-	0/20/34/34	0/1/3/3
2	DTP	B	301	-	-	0/20/34/34	0/1/3/3
2	DTP	C	301	-	-	0/20/34/34	0/1/3/3
2	DTP	D	301	-	-	0/20/34/34	0/1/3/3
2	DTP	E	301	-	-	0/20/34/34	0/1/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DTP	F	301	-	-	0/20/34/34	0/1/3/3
2	DTP	G	301	-	-	0/20/34/34	0/1/3/3
2	DTP	H	301	-	-	0/20/34/34	0/1/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	301	DTP	C8-N7	-2.47	1.29	1.34
2	D	301	DTP	C8-N7	-2.33	1.30	1.34
2	C	301	DTP	C8-N7	-2.20	1.30	1.34
2	E	301	DTP	C8-N7	-2.12	1.30	1.34
2	H	301	DTP	C8-N7	-2.12	1.30	1.34
2	D	301	DTP	PB-O3A	-2.04	1.56	1.59
2	G	301	DTP	C8-N7	-2.02	1.30	1.34

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	DTP	N3-C2-N1	-9.72	120.58	128.71
2	G	301	DTP	N3-C2-N1	-9.47	120.79	128.71
2	H	301	DTP	N3-C2-N1	-9.39	120.86	128.71
2	D	301	DTP	N3-C2-N1	-9.35	120.89	128.71
2	C	301	DTP	N3-C2-N1	-9.24	120.98	128.71
2	B	301	DTP	N3-C2-N1	-9.15	121.06	128.71
2	A	301	DTP	N3-C2-N1	-8.80	121.35	128.71
2	E	301	DTP	N3-C2-N1	-8.79	121.36	128.71
2	C	301	DTP	N3-C4-N9	5.11	134.66	125.43
2	F	301	DTP	N3-C4-N9	5.06	134.57	125.43
2	D	301	DTP	N3-C4-N9	4.99	134.44	125.43
2	E	301	DTP	N3-C4-N9	4.99	134.43	125.43
2	H	301	DTP	N3-C4-N9	4.87	134.22	125.43
2	G	301	DTP	N3-C4-N9	4.86	134.20	125.43
2	A	301	DTP	N3-C4-N9	4.63	133.80	125.43
2	B	301	DTP	N3-C4-N9	4.57	133.68	125.43
2	E	301	DTP	C5'-C4'-C3'	-4.03	89.54	114.76
2	G	301	DTP	C5'-C4'-C3'	-3.92	90.26	114.76
2	B	301	DTP	C5'-C4'-C3'	-3.90	90.35	114.76
2	F	301	DTP	C5'-C4'-C3'	-3.86	90.60	114.76
2	A	301	DTP	C5'-C4'-C3'	-3.85	90.69	114.76
2	D	301	DTP	C5'-C4'-C3'	-3.84	90.72	114.76
2	C	301	DTP	C5'-C4'-C3'	-3.77	91.19	114.76
2	H	301	DTP	C5'-C4'-C3'	-3.75	91.28	114.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	DTP	O5'-C5'-C4'	3.02	120.03	108.94
2	G	301	DTP	O5'-C5'-C4'	2.92	119.64	108.94
2	A	301	DTP	O5'-C5'-C4'	2.91	119.62	108.94
2	C	301	DTP	O5'-C5'-C4'	2.87	119.47	108.94
2	B	301	DTP	O5'-C5'-C4'	2.86	119.44	108.94
2	H	301	DTP	O5'-C5'-C4'	2.84	119.37	108.94
2	H	301	DTP	C5-C4-N3	-2.78	119.64	125.70
2	D	301	DTP	C5-C4-N3	-2.78	119.65	125.70
2	E	301	DTP	O5'-C5'-C4'	2.77	119.10	108.94
2	D	301	DTP	O5'-C5'-C4'	2.77	119.09	108.94
2	E	301	DTP	C5-C4-N3	-2.76	119.69	125.70
2	C	301	DTP	C5-C4-N3	-2.71	119.80	125.70
2	G	301	DTP	C5-C4-N3	-2.71	119.80	125.70
2	H	301	DTP	C2-N3-C4	2.70	121.69	114.01
2	F	301	DTP	C2-N3-C4	2.65	121.54	114.01
2	F	301	DTP	C5-C4-N3	-2.62	119.99	125.70
2	D	301	DTP	C2-N3-C4	2.62	121.46	114.01
2	B	301	DTP	C5-C4-N3	-2.61	120.03	125.70
2	C	301	DTP	C2-N3-C4	2.53	121.21	114.01
2	A	301	DTP	C5-C4-N3	-2.52	120.20	125.70
2	G	301	DTP	C2-N3-C4	2.52	121.19	114.01
2	B	301	DTP	C2-N3-C4	2.51	121.17	114.01
2	E	301	DTP	C2-N3-C4	2.48	121.07	114.01
2	G	301	DTP	C4-C5-N7	-2.43	107.44	109.52
2	A	301	DTP	C2-N3-C4	2.35	120.71	114.01
2	B	301	DTP	C4-C5-N7	-2.27	107.58	109.52
2	C	301	DTP	C4-C5-N7	-2.22	107.62	109.52
2	D	301	DTP	C4-C5-N7	-2.17	107.66	109.52
2	E	301	DTP	C4-C5-N7	-2.15	107.68	109.52
2	E	301	DTP	O4'-C1'-N9	2.14	111.71	107.68
2	C	301	DTP	O3'-C3'-C2'	2.12	118.78	110.78
2	F	301	DTP	O4'-C1'-N9	2.11	111.65	107.68
2	A	301	DTP	O2B-PB-O3B	2.04	114.82	105.14
2	A	301	DTP	O3'-C3'-C2'	2.02	118.39	110.78
2	E	301	DTP	O3'-C3'-C2'	2.02	118.39	110.78

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	229/241 (95%)	0.17	18 (7%) 13 11	17, 39, 92, 113	0
1	B	229/241 (95%)	0.09	9 (3%) 37 37	16, 41, 85, 104	0
1	C	229/241 (95%)	0.08	12 (5%) 26 26	12, 31, 81, 95	0
1	D	229/241 (95%)	-0.03	7 (3%) 47 47	10, 29, 70, 103	0
1	E	229/241 (95%)	0.19	16 (6%) 16 14	19, 44, 91, 109	0
1	F	229/241 (95%)	0.22	16 (6%) 16 14	16, 36, 76, 120	0
1	G	229/241 (95%)	0.10	14 (6%) 21 20	19, 39, 75, 107	0
1	H	229/241 (95%)	0.30	20 (8%) 10 9	20, 46, 90, 107	0
All	All	1832/1928 (95%)	0.14	112 (6%) 21 20	10, 38, 87, 120	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	72	VAL	7.1
1	H	134	ARG	6.8
1	F	73	ALA	6.1
1	F	75	TRP	6.1
1	H	131	LEU	6.0
1	C	130	LEU	5.6
1	F	128	GLU	5.1
1	A	130	LEU	4.9
1	A	75	TRP	4.7
1	F	132	GLN	4.7
1	A	134	ARG	4.6
1	D	244	ASN	4.5
1	G	75	TRP	4.5
1	F	129	LYS	4.4
1	H	132	GLN	4.1
1	G	76	GLN	4.0

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Mol	Chain	Res	Type	RSRZ
1	H	133	ALA	3.9
1	H	37	GLY	3.8
1	C	129	LYS	3.8
1	H	72	VAL	3.7
1	A	132	GLN	3.7
1	B	134	ARG	3.6
1	B	75	TRP	3.6
1	G	72	VAL	3.5
1	E	79	GLN	3.5
1	F	77	ASN	3.5
1	E	256	PHE	3.5
1	H	79	GLN	3.4
1	F	37	GLY	3.4
1	H	128	GLU	3.4
1	A	78	ILE	3.3
1	C	69	THR	3.3
1	E	240	GLU	3.3
1	E	131	LEU	3.3
1	C	134	ARG	3.3
1	A	72	VAL	3.3
1	A	212	LYS	3.1
1	E	75	TRP	3.1
1	A	77	ASN	3.1
1	A	261	THR	3.0
1	F	131	LEU	3.0
1	E	132	GLN	3.0
1	G	132	GLN	3.0
1	G	261	THR	3.0
1	F	74	THR	3.0
1	E	78	ILE	2.9
1	H	130	LEU	2.9
1	A	76	GLN	2.9
1	E	76	GLN	2.9
1	A	92	LEU	2.9
1	F	126	PHE	2.9
1	F	78	ILE	2.9
1	H	78	ILE	2.9
1	B	79	GLN	2.8
1	A	93	GLY	2.8
1	E	134	ARG	2.8
1	E	72	VAL	2.8
1	H	75	TRP	2.7

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Mol	Chain	Res	Type	RSRZ
1	H	242	LEU	2.7
1	C	64	GLU	2.7
1	C	79	GLN	2.7
1	G	74	THR	2.7
1	G	71	PRO	2.6
1	B	128	GLU	2.6
1	F	204	TYR	2.6
1	H	268	ARG	2.6
1	H	244	ASN	2.6
1	E	135	LYS	2.6
1	B	131	LEU	2.5
1	A	129	LYS	2.5
1	H	277	LEU	2.5
1	D	204	TYR	2.5
1	A	69	THR	2.5
1	G	69	THR	2.5
1	G	79	GLN	2.5
1	C	132	GLN	2.4
1	F	76	GLN	2.4
1	C	135	LYS	2.4
1	D	134	ARG	2.4
1	A	79	GLN	2.4
1	E	260	VAL	2.4
1	H	65	TRP	2.4
1	H	135	LYS	2.3
1	G	70	GLU	2.3
1	B	135	LYS	2.3
1	F	69	THR	2.3
1	D	76	GLN	2.3
1	A	259	GLU	2.2
1	E	258	GLU	2.2
1	A	37	GLY	2.2
1	E	259	GLU	2.2
1	C	261	THR	2.2
1	A	73	ALA	2.2
1	E	128	GLU	2.2
1	C	259	GLU	2.2
1	D	77	ASN	2.2
1	F	178	TRP	2.1
1	H	129	LYS	2.1
1	B	130	LEU	2.1
1	B	277	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	59	THR	2.1
1	G	268	ARG	2.1
1	D	261	THR	2.1
1	G	128	GLU	2.1
1	B	243	MET	2.0
1	H	261	THR	2.0
1	C	75	TRP	2.0
1	H	204	TYR	2.0
1	D	75	TRP	2.0
1	G	259	GLU	2.0
1	C	131	LEU	2.0
1	G	258	GLU	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, 95th 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 ²)	Q<0.9
2	DTP	D	301	30/30	0.19	0.26	22,35,54,55	0
2	DTP	C	301	30/30	0.18	-0.03	22,35,54,56	0
2	DTP	B	301	30/30	0.18	-0.17	21,36,55,55	0
2	DTP	E	301	30/30	0.18	-0.32	22,36,54,56	0
2	DTP	F	301	30/30	0.17	-0.34	22,35,54,55	0
2	DTP	H	301	30/30	0.16	-0.34	22,36,54,55	0
2	DTP	A	301	30/30	0.17	-0.38	22,35,55,56	0
2	DTP	G	301	30/30	0.14	-0.63	22,36,55,56	0

6.5 Other polymers ⓘ

There are no such residues in this entry.