



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:48 AM GMT

PDB ID : 3KRP
Title : Mint heterotetrameric geranyl pyrophosphate synthase in complex with magnesium and GPP
Authors : Chang, T.-H.; Hsieh, F.-L.; Ko, T.-P.; Wang, A.H.-J.
Deposited on : 2009-11-19
Resolution : 2.42 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>

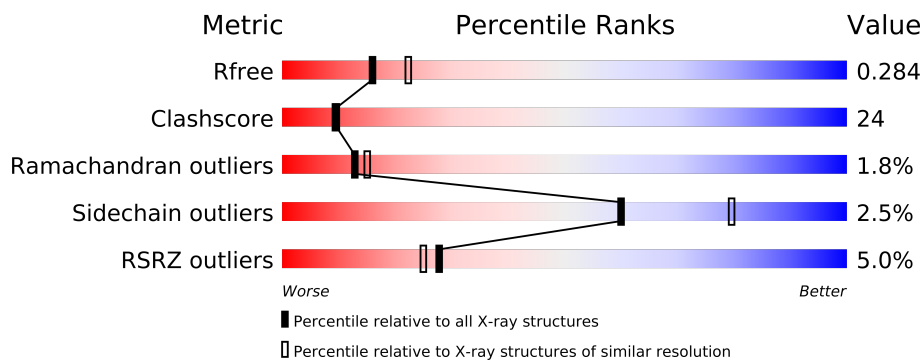
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.42 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	2496 (2.44-2.40)
Clashscore	79885	3124 (2.44-2.40)
Ramachandran outliers	78287	3067 (2.44-2.40)
Sidechain outliers	78261	3068 (2.44-2.40)
RSRZ outliers	66119	2499 (2.44-2.40)

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	295	
1	D	295	
2	B	274	
2	C	274	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9010 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 Geranyl diphosphate synthase large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	S	0	0	0
			2120	1338	372	392	18			
1	D	284	Total	C	N	O	S	0	0	0
			2155	1359	378	400	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q9SBR3
D	1	MET	-	EXPRESSION TAG	UNP Q9SBR3

- Molecule 2 is a protein called Geranyl diphosphate synthase small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	256	Total	C	N	O	S	0	0	0
			1929	1216	334	365	14			
2	C	259	Total	C	N	O	S	0	0	0
			1954	1233	339	368	14			

There are 18 discrepancies between the modelled and reference sequences:

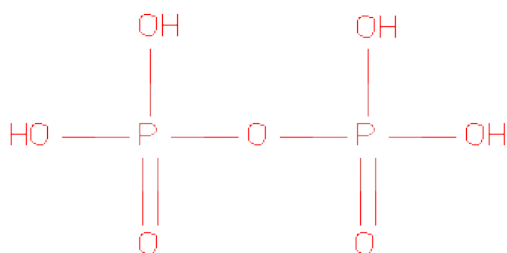
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	EXPRESSION TAG	UNP Q9SBR4
B	267	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	268	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	269	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	270	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	271	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	272	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	273	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	274	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	1	MET	-	EXPRESSION TAG	UNP Q9SBR4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	267	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	268	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	269	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	270	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	271	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	272	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	273	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	274	HIS	-	EXPRESSION TAG	UNP Q9SBR4

- Molecule 3 is PYROPHOSPHATE (three-letter code: PPV) (formula: $\text{H}_4\text{O}_7\text{P}_2$).

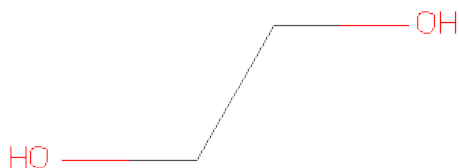


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 9 7 2	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

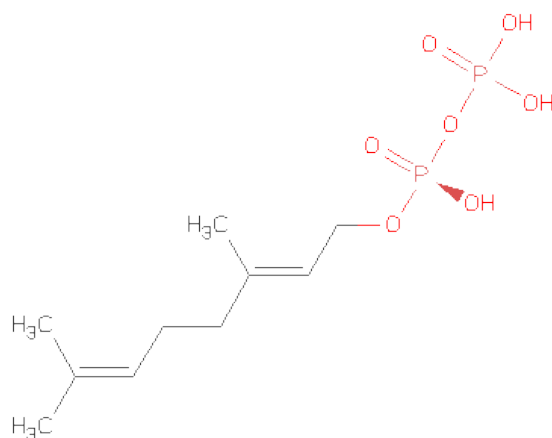
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Mg 2 2	0	0
4	D	2	Total Mg 2 2	0	0

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			4	2	2		

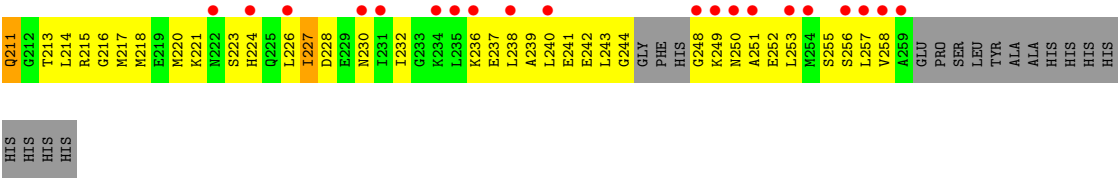
- Molecule 6 is GERANYL DIPHOSPHATE (three-letter code: GPP) (formula: $C_{10}H_{20}O_7P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	D	1	Total	C	O	P	0	0
			19	10	7	2		

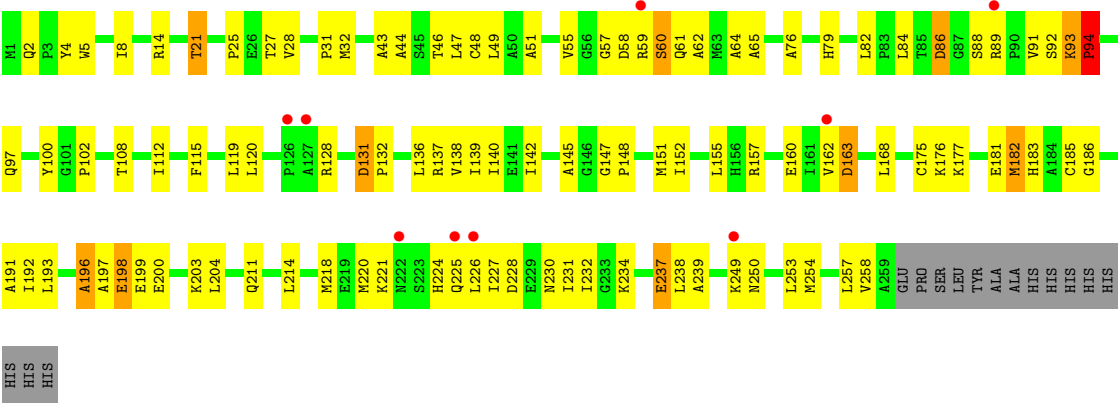
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	224	Total 224	O 224	0	0
7	B	177	Total 177	O 177	0	0
7	C	221	Total 221	O 221	0	0
7	D	194	Total 194	O 194	0	0



● Molecule 2: Geranyl diphosphate synthase small subunit

Chain C:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.26Å 109.02Å 182.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.42 29.62 – 2.24	Depositor EDS
% Data completeness (in resolution range)	85.9 (30.00-2.42) 83.1 (29.62-2.24)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.82 (at 2.24Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.223 , 0.284 0.225 , 0.284	Depositor DCC
R_{free} test set	1790 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	1.003	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 43365 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9010	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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: MG, GPP, EDO, PPV

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	0/2155	0.59	0/2903
1	D	0.37	0/2190	0.59	0/2948
2	B	0.35	0/1965	0.57	0/2656
2	C	0.37	0/1993	0.57	0/2695
All	All	0.38	0/8303	0.58	0/11202

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	2120	0	2146	81	0
1	D	2155	0	2182	130	0
2	B	1929	0	1928	117	0
2	C	1954	0	1948	93	0
3	A	9	0	0	0	0
4	A	2	0	0	0	0
4	D	2	0	0	0	0
5	C	4	0	6	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	19	0	17	2	0
7	A	224	0	0	13	0
7	B	177	0	0	16	0
7	C	221	0	0	11	0
7	D	194	0	0	13	0
All	All	9010	0	8227	401	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 24.

All (401) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:23:ARG:HB2	2:B:24:PRO:HD2	1.37	1.06
2:B:249:LYS:HE2	2:B:253:LEU:HD13	1.46	0.98
2:B:8:ILE:HG12	2:B:59:ARG:HH12	1.28	0.97
2:C:163:ASP:HA	7:C:702:HOH:O	1.65	0.95
1:A:168:VAL:HG13	1:A:172:HIS:HB3	1.48	0.94
2:B:249:LYS:HD3	2:B:253:LEU:HB2	1.49	0.93
2:C:221:LYS:HD3	2:C:227:ILE:HG22	1.50	0.90
2:C:43:ALA:O	2:C:47:LEU:HB2	1.70	0.90
2:C:14:ARG:HG2	7:C:535:HOH:O	1.73	0.88
1:A:139:VAL:HG13	1:A:191:LEU:HD22	1.57	0.85
2:B:248:GLY:N	2:B:251:ALA:HB3	1.91	0.84
1:D:260:PHE:HD1	1:D:263:ARG:HH21	1.25	0.84
2:B:55:VAL:HG11	2:B:196:ALA:HB2	1.60	0.82
2:C:8:ILE:HG23	2:C:49:LEU:HD12	1.63	0.81
2:C:28:VAL:HG22	1:D:157:VAL:CG2	2.11	0.81
2:B:2:GLN:HB2	2:B:3:PRO:HD3	1.63	0.80
1:A:289:TYR:O	1:A:293:ARG:HG2	1.82	0.79
1:D:135:ARG:HG3	1:D:135:ARG:HH11	1.46	0.79
1:A:277:PRO:HD2	1:A:278:HIS:CE1	2.17	0.78
1:D:10:ARG:HD2	7:D:394:HOH:O	1.84	0.78
1:D:139:VAL:HG13	1:D:191:LEU:HD22	1.64	0.77
1:A:76:HIS:CE1	1:A:80:LEU:HD11	2.20	0.77
2:C:214:LEU:O	2:C:218:MET:HG2	1.86	0.76
1:D:9:LEU:O	1:D:13:LYS:HD3	1.86	0.75
2:C:199:GLU:O	2:C:203:LYS:HG2	1.87	0.74
1:A:256:LYS:HE2	7:A:378:HOH:O	1.87	0.73
1:D:216:LEU:O	1:D:220:VAL:HG23	1.88	0.73
2:B:2:GLN:HB2	2:B:3:PRO:CD	2.19	0.72
1:A:183:ALA:HB3	7:A:341:HOH:O	1.89	0.71
2:C:97:GLN:NE2	7:C:773:HOH:O	2.22	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:47:ARG:HB2	1:A:48:PRO:HD3	1.73	0.71
1:A:1:MET:N	7:A:352:HOH:O	2.23	0.71
2:B:138:VAL:O	2:B:142:ILE:HG13	1.91	0.71
2:B:249:LYS:O	2:B:253:LEU:N	2.24	0.70
2:C:137:ARG:HD3	2:C:192:ILE:HD13	1.72	0.70
2:B:28:VAL:O	2:B:31:PRO:HD2	1.91	0.70
2:C:230:ASN:O	2:C:234:LYS:HG3	1.92	0.70
1:A:165:MET:HE1	7:A:757:HOH:O	1.91	0.70
2:C:182:MET:HE3	2:C:182:MET:HA	1.73	0.70
2:B:221:LYS:HG3	2:B:227:ILE:HG21	1.72	0.70
2:C:221:LYS:HD3	2:C:227:ILE:CG2	2.20	0.69
1:A:256:LYS:HE3	7:A:379:HOH:O	1.91	0.69
2:B:248:GLY:O	2:B:252:GLU:HB2	1.91	0.69
1:A:286:LEU:O	1:A:290:ILE:HG13	1.93	0.69
2:B:23:ARG:HB2	2:B:24:PRO:CD	2.19	0.69
2:B:137:ARG:HD3	2:B:192:ILE:HD13	1.75	0.69
2:B:253:LEU:HD12	2:B:256:SER:OG	1.93	0.68
1:D:20:GLU:HA	1:D:40:LEU:HD11	1.75	0.68
2:B:1:MET:HG3	2:B:2:GLN:H	1.57	0.68
1:D:135:ARG:NH1	7:D:368:HOH:O	2.27	0.68
2:B:25:PRO:HB2	2:B:27:THR:HG22	1.76	0.68
2:C:62:ALA:HA	2:C:193:LEU:HD13	1.74	0.67
1:A:19:LEU:HD22	1:A:39:LEU:HD21	1.75	0.67
1:D:168:VAL:HG23	1:D:251:LEU:HD21	1.75	0.67
1:A:64:THR:HG23	1:A:129:LYS:O	1.94	0.67
1:A:149:SER:HB2	7:A:493:HOH:O	1.92	0.67
1:D:252:ILE:O	1:D:256:LYS:HB2	1.94	0.67
2:B:227:ILE:O	2:B:227:ILE:HG13	1.93	0.67
2:B:220:MET:HB3	2:B:224:HIS:HE1	1.59	0.67
2:B:8:ILE:HG12	2:B:59:ARG:NH1	2.06	0.67
2:C:58:ASP:OD1	2:C:61:GLN:HG3	1.95	0.66
2:B:168:LEU:HD21	2:B:226:LEU:HB2	1.77	0.66
2:C:86:ASP:OD2	1:D:106:GLU:HG3	1.96	0.66
1:D:260:PHE:HD1	1:D:263:ARG:NH2	1.93	0.66
2:B:177:LYS:O	7:B:508:HOH:O	2.13	0.66
1:D:208:LYS:HD2	7:D:694:HOH:O	1.95	0.66
2:B:249:LYS:CD	2:B:253:LEU:HB2	2.23	0.65
1:D:170:LEU:O	1:D:174:GLU:HB2	1.96	0.65
1:A:96:GLY:C	1:A:97:LYS:HD2	2.15	0.65
1:D:80:LEU:HD23	6:D:900:GPP:H92	1.77	0.65
1:A:167:GLU:HG2	7:A:546:HOH:O	1.96	0.65
1:A:90:ASN:O	7:A:412:HOH:O	2.13	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:43:ALA:HB3	2:B:183:HIS:HE1	1.62	0.65
2:B:43:ALA:HB3	2:B:183:HIS:CE1	2.32	0.64
1:D:151:GLY:O	1:D:179:HIS:HB3	1.97	0.64
2:B:188:ALA:O	2:B:192:ILE:HG13	1.97	0.64
2:B:92:SER:C	2:B:94:PRO:HD2	2.18	0.64
1:A:245:LYS:O	1:A:245:LYS:HD2	1.97	0.64
1:D:208:LYS:HE2	7:D:529:HOH:O	1.96	0.64
1:A:276:HIS:HB3	1:A:278:HIS:ND1	2.12	0.64
1:D:251:LEU:HG	7:D:665:HOH:O	1.96	0.64
2:B:177:LYS:NZ	2:B:181:GLU:OE2	2.32	0.63
1:A:153:VAL:O	1:A:157:VAL:HG23	1.97	0.63
2:B:214:LEU:O	2:B:218:MET:HG2	1.98	0.63
2:C:214:LEU:HD21	2:C:232:ILE:HG23	1.81	0.63
2:B:236:LYS:O	2:B:240:LEU:HG	1.97	0.63
1:A:149:SER:CB	7:A:493:HOH:O	2.47	0.63
2:C:175:CYS:SG	2:C:220:MET:HG3	2.38	0.63
2:C:191:ALA:HB2	2:C:204:LEU:HD12	1.80	0.63
2:C:64:ALA:HB3	7:C:350:HOH:O	1.98	0.62
2:B:136:LEU:O	2:B:136:LEU:HD12	1.99	0.62
2:B:71:LEU:HD22	2:B:112:ILE:HG23	1.82	0.62
2:C:136:LEU:O	2:C:140:ILE:HG13	2.00	0.62
2:B:1:MET:HG3	2:B:2:GLN:N	2.14	0.62
2:C:28:VAL:HG22	1:D:157:VAL:HG21	1.82	0.62
1:D:187:GLY:O	1:D:191:LEU:HG	1.99	0.62
1:A:30:LYS:HE2	1:A:108:VAL:HG21	1.81	0.62
2:C:51:ALA:HA	2:C:254:MET:HE3	1.82	0.62
1:D:132:PRO:HB2	1:D:134:GLU:OE2	2.00	0.61
1:D:220:VAL:HG22	7:D:576:HOH:O	2.01	0.61
1:A:86:PRO:HG2	1:A:106:GLU:OE2	2.01	0.61
1:D:139:VAL:CG1	1:D:191:LEU:HD22	2.31	0.61
1:D:93:LEU:HG	1:D:234:GLY:O	2.00	0.61
2:C:65:ALA:HA	2:C:120:LEU:HD21	1.83	0.60
2:C:25:PRO:HB2	2:C:27:THR:HG22	1.83	0.60
2:C:227:ILE:O	2:C:227:ILE:HG23	2.01	0.60
1:D:93:LEU:HD11	1:D:96:GLY:HA2	1.83	0.60
1:A:270:GLU:HA	1:A:273:LEU:HG	1.84	0.60
1:A:76:HIS:NE2	1:A:80:LEU:HD11	2.17	0.60
1:A:61:ASP:OD2	1:A:63:SER:HB3	2.03	0.59
2:B:179:TYR:OH	7:B:306:HOH:O	2.16	0.59
1:D:30:LYS:HE2	1:D:108:VAL:HG21	1.84	0.59
1:D:7:TYR:HA	1:D:10:ARG:HH11	1.67	0.59
2:B:221:LYS:HG3	2:B:227:ILE:CG2	2.32	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:237:GLU:O	2:C:239:ALA:N	2.36	0.59
2:B:31:PRO:HA	2:B:100:TYR:CE2	2.38	0.58
2:C:162:VAL:HG12	2:C:163:ASP:OD1	2.04	0.58
1:D:20:GLU:HA	1:D:40:LEU:CD1	2.33	0.58
1:D:256:LYS:HA	1:D:259:GLU:CD	2.24	0.58
2:C:181:GLU:HA	2:C:181:GLU:OE2	2.04	0.58
2:B:240:LEU:HD21	2:B:258:VAL:HG21	1.86	0.58
2:B:257:LEU:HD21	7:B:505:HOH:O	2.03	0.58
1:D:177:HIS:ND1	1:D:218:GLN:HG2	2.19	0.57
1:A:228:LYS:NZ	1:A:228:LYS:HB2	2.19	0.57
2:C:92:SER:C	2:C:94:PRO:HD2	2.25	0.57
2:C:93:LYS:N	2:C:94:PRO:HD2	2.20	0.57
2:B:147:GLY:HA3	7:B:296:HOH:O	2.04	0.57
2:C:79:HIS:HA	2:C:82:LEU:HG	1.87	0.57
1:D:83:ASP:O	1:D:89:ASP:HB2	2.05	0.57
2:B:137:ARG:NH1	7:B:499:HOH:O	2.37	0.57
2:B:23:ARG:CB	2:B:24:PRO:HD2	2.24	0.56
2:B:213:THR:O	2:B:217:MET:HB2	2.05	0.56
1:D:222:ASP:HB3	1:D:249:PRO:HD2	1.87	0.56
1:D:208:LYS:HG2	7:D:529:HOH:O	2.05	0.56
1:D:281:ALA:HB3	1:D:282:PRO:HD3	1.87	0.56
2:B:120:LEU:HD11	2:B:142:ILE:HD12	1.87	0.56
1:A:88:MET:HE3	1:A:160:VAL:HG21	1.87	0.56
2:C:58:ASP:HB2	7:C:634:HOH:O	2.05	0.56
2:B:54:LEU:HB2	2:B:250:ASN:OD1	2.05	0.56
1:A:133:PRO:O	1:A:137:VAL:HG23	2.05	0.56
1:D:180:LYS:HG2	6:D:900:GPP:H62	1.86	0.56
2:C:102:PRO:HB2	1:D:87:CYS:HB2	1.87	0.56
2:B:55:VAL:CG1	2:B:196:ALA:HB2	2.33	0.56
1:D:252:ILE:O	1:D:256:LYS:HE3	2.06	0.56
1:D:7:TYR:HA	1:D:10:ARG:NH1	2.21	0.55
1:D:263:ARG:HB2	1:D:263:ARG:NH1	2.21	0.55
2:B:214:LEU:HD23	2:B:258:VAL:HG12	1.88	0.55
2:B:9:GLU:OE1	7:B:503:HOH:O	2.18	0.55
2:C:43:ALA:HB3	2:C:183:HIS:HE1	1.72	0.55
2:B:220:MET:HB3	2:B:224:HIS:CE1	2.40	0.55
2:B:168:LEU:HD21	2:B:226:LEU:CB	2.37	0.55
1:A:205:LYS:HD3	1:A:271:GLN:O	2.07	0.55
2:C:221:LYS:HD2	2:C:225:GLN:HA	1.89	0.55
2:C:4:TYR:CE2	2:C:8:ILE:HD11	2.42	0.55
1:A:75:ILE:O	1:A:78:MET:HG3	2.07	0.55
1:D:289:TYR:O	1:D:293:ARG:HB3	2.07	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:44:ALA:CB	2:C:183:HIS:HA	2.37	0.55
1:A:95:ARG:HG2	1:A:95:ARG:NH1	2.22	0.54
1:D:68:ALA:HA	1:D:124:VAL:HG22	1.89	0.54
1:A:87:CYS:HB2	2:B:102:PRO:HB2	1.89	0.54
2:B:137:ARG:HD2	7:B:499:HOH:O	2.07	0.54
1:A:95:ARG:HH11	1:A:95:ARG:HG2	1.73	0.54
1:D:294:ASP:O	1:D:295:ASN:HB3	2.07	0.54
1:D:229:SER:OG	1:D:232:GLU:HG3	2.08	0.53
1:D:280:ALA:O	1:D:284:ILE:HG13	2.08	0.53
1:D:37:TYR:O	1:D:41:ALA:HB2	2.08	0.53
1:A:140:LEU:HD13	2:B:139:ILE:HD13	1.91	0.53
2:B:93:LYS:N	2:B:94:PRO:HD2	2.22	0.53
2:C:92:SER:CB	2:C:94:PRO:HD2	2.38	0.53
1:A:31:ILE:HD11	2:B:156:HIS:CD2	2.43	0.53
2:C:108:THR:O	2:C:112:ILE:HG12	2.08	0.53
1:D:177:HIS:ND1	1:D:218:GLN:CG	2.72	0.53
1:A:281:ALA:HB3	1:A:282:PRO:HD3	1.91	0.53
1:D:149:SER:O	1:D:154:ALA:HB2	2.08	0.53
2:B:91:VAL:O	2:B:91:VAL:HG23	2.09	0.53
1:D:123:HIS:O	1:D:127:ALA:HB3	2.09	0.53
2:B:155:LEU:HD23	2:B:178:LYS:HE3	1.91	0.53
1:D:223:ILE:O	1:D:227:THR:HG23	2.10	0.52
1:D:66:MET:HB3	1:D:67:PRO:HD3	1.92	0.52
2:C:84:LEU:HD12	2:C:89:ARG:HG2	1.92	0.52
1:D:135:ARG:HG3	1:D:135:ARG:NH1	2.21	0.52
1:A:186:GLN:NE2	1:A:211:ASN:OD1	2.42	0.52
1:D:75:ILE:HD13	1:D:147:ILE:HD13	1.90	0.52
2:B:41:ALA:HB1	7:B:300:HOH:O	2.10	0.52
2:C:214:LEU:HD13	2:C:214:LEU:C	2.30	0.52
2:B:62:ALA:HA	2:B:193:LEU:HD13	1.91	0.52
2:B:258:VAL:HG22	7:B:616:HOH:O	2.11	0.51
1:A:94:ARG:HG2	1:A:95:ARG:HG3	1.93	0.51
2:C:157:ARG:O	2:C:160:GLU:HB2	2.10	0.51
1:D:178:HIS:O	1:D:182:ALA:HB3	2.11	0.51
2:C:43:ALA:HB3	2:C:183:HIS:CE1	2.46	0.51
1:A:89:ASP:OD1	7:A:733:HOH:O	2.18	0.51
2:C:8:ILE:CG2	2:C:49:LEU:HD12	2.37	0.51
1:D:56:GLU:HA	1:D:60:GLY:O	2.11	0.50
1:D:215:LEU:O	1:D:219:VAL:HG23	2.10	0.50
1:D:222:ASP:CG	1:D:247:THR:HB	2.30	0.50
1:D:279:ARG:O	1:D:282:PRO:HD2	2.10	0.50
2:B:17:LYS:HG3	2:B:37:PHE:CE2	2.46	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:4:TYR:O	2:B:8:ILE:HG13	2.11	0.50
2:B:249:LYS:HD2	2:B:253:LEU:HD22	1.93	0.50
2:C:51:ALA:HA	2:C:254:MET:CE	2.41	0.50
1:D:93:LEU:HD11	1:D:96:GLY:CA	2.41	0.50
2:B:5:TRP:HZ2	2:B:253:LEU:HG	1.77	0.50
2:B:146:GLY:O	2:B:152:ILE:HG12	2.11	0.50
2:C:177:LYS:HA	2:C:181:GLU:HG2	1.93	0.50
2:C:88:SER:OG	1:D:107:SER:HB3	2.12	0.50
1:D:250:LYS:HD3	7:D:665:HOH:O	2.10	0.49
1:D:183:ALA:HB3	7:D:313:HOH:O	2.12	0.49
2:B:242:GLU:O	2:B:243:LEU:HD23	2.12	0.49
1:A:148:GLY:O	1:A:153:VAL:HB	2.12	0.49
2:B:75:ALA:CB	2:B:113:VAL:HG23	2.42	0.49
2:C:211:GLN:HG2	2:C:258:VAL:HG12	1.95	0.49
1:D:248:TYR:HB2	1:D:249:PRO:HD3	1.95	0.49
2:B:162:VAL:O	2:B:162:VAL:HG12	2.13	0.49
1:A:209:PHE:CE1	1:A:272:LEU:HD11	2.48	0.49
2:C:221:LYS:CD	2:C:227:ILE:HG22	2.33	0.49
1:D:212:CYS:C	1:D:214:GLY:H	2.14	0.49
1:D:288:ASN:O	1:D:292:TYR:HD1	1.96	0.49
2:B:221:LYS:HD3	2:B:227:ILE:HB	1.95	0.49
2:B:239:ALA:O	2:B:242:GLU:HB2	2.13	0.49
1:D:69:ALA:O	1:D:72:VAL:HG22	2.12	0.49
2:B:101:GLY:O	2:B:105:GLU:HG3	2.13	0.48
2:B:179:TYR:CD1	2:B:216:GLY:HA3	2.48	0.48
1:A:8:MET:SD	1:A:49:MET:HG3	2.53	0.48
2:C:59:ARG:HG3	2:C:60:SER:N	2.28	0.48
2:C:8:ILE:HG22	2:C:46:THR:HG22	1.95	0.48
2:C:93:LYS:N	2:C:94:PRO:CD	2.76	0.48
1:A:148:GLY:HA3	7:B:335:HOH:O	2.11	0.48
2:B:230:ASN:HB3	7:B:279:HOH:O	2.13	0.48
1:A:208:LYS:HE3	1:A:271:GLN:CD	2.34	0.48
2:B:21:THR:HG22	2:B:21:THR:O	2.13	0.48
2:B:71:LEU:HD12	2:B:116:GLY:HA2	1.95	0.48
1:D:288:ASN:ND2	7:D:310:HOH:O	2.47	0.48
1:A:152:LEU:HA	1:A:179:HIS:O	2.12	0.48
2:C:176:LYS:HE3	7:C:286:HOH:O	2.12	0.48
2:C:55:VAL:HG21	2:C:196:ALA:HB2	1.95	0.48
1:A:247:THR:HG21	7:A:736:HOH:O	2.13	0.48
1:D:47:ARG:HG2	1:D:185:LEU:HD23	1.95	0.48
7:C:291:HOH:O	1:D:149:SER:HB2	2.12	0.48
1:A:70:CYS:O	1:A:74:MET:HG3	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:253:LEU:C	2:B:255:SER:N	2.67	0.47
2:C:197:ALA:O	2:C:198:GLU:C	2.52	0.47
1:A:168:VAL:CG1	1:A:172:HIS:HB3	2.33	0.47
2:C:168:LEU:HD13	2:C:231:ILE:HG13	1.95	0.47
2:B:20:ILE:HD12	2:B:36:THR:HG21	1.95	0.47
1:D:78:MET:HE3	1:D:113:GLY:O	2.14	0.47
1:D:4:PHE:HB2	1:D:8:MET:HE1	1.96	0.47
1:A:19:LEU:HB3	1:A:40:LEU:HD21	1.96	0.47
2:C:249:LYS:O	2:C:253:LEU:HG	2.15	0.47
2:B:173:TYR:CE2	1:D:29:LEU:HD21	2.49	0.47
2:C:228:ASP:OD1	2:C:230:ASN:HB2	2.15	0.47
2:B:228:ASP:OD2	2:B:230:ASN:HB2	2.15	0.47
2:B:258:VAL:HG13	7:B:616:HOH:O	2.14	0.47
2:B:203:LYS:HD2	2:B:244:GLY:H	1.79	0.47
2:C:139:ILE:HG21	1:D:140:LEU:HD13	1.96	0.47
1:D:164:GLY:O	1:D:166:ALA:N	2.48	0.47
1:A:89:ASP:HB3	1:A:91:ASP:OD1	2.14	0.47
2:B:248:GLY:N	7:B:714:HOH:O	2.48	0.47
2:C:48:CYS:HA	2:C:186:GLY:O	2.15	0.47
2:C:131:ASP:N	2:C:132:PRO:HD2	2.30	0.47
2:B:1:MET:CG	2:B:2:GLN:N	2.77	0.47
2:C:128:ARG:HB3	2:C:128:ARG:NH1	2.30	0.47
1:D:4:PHE:HB2	1:D:8:MET:CE	2.45	0.46
1:D:222:ASP:OD2	1:D:247:THR:HB	2.15	0.46
1:D:27:GLU:HB2	7:D:578:HOH:O	2.14	0.46
2:B:197:ALA:O	2:B:201:ILE:HG13	2.14	0.46
2:B:216:GLY:O	2:B:220:MET:HG2	2.15	0.46
2:B:226:LEU:O	2:B:228:ASP:N	2.47	0.46
1:D:181:THR:HG21	1:D:218:GLN:HB2	1.97	0.46
2:B:1:MET:O	2:B:5:TRP:CD1	2.68	0.46
2:C:46:THR:HG23	5:C:902:EDO:O2	2.16	0.46
1:D:93:LEU:HD12	1:D:94:ARG:H	1.81	0.46
1:D:214:GLY:C	1:D:216:LEU:N	2.68	0.46
1:A:19:LEU:HD22	1:A:39:LEU:CD2	2.45	0.46
2:C:93:LYS:O	2:C:94:PRO:C	2.54	0.46
1:D:223:ILE:HA	1:D:226:VAL:HG12	1.97	0.46
2:C:32:MET:HA	2:C:108:THR:CG2	2.46	0.46
1:D:286:LEU:O	1:D:290:ILE:HG13	2.15	0.46
1:D:157:VAL:HG23	1:D:158:VAL:N	2.31	0.46
1:A:25:MET:O	1:A:26:LYS:HB3	2.16	0.46
2:C:151:MET:O	2:C:155:LEU:HG	2.16	0.46
2:C:28:VAL:HG22	1:D:157:VAL:HG23	1.93	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:186:GLN:HG3	1:A:207:ARG:HG3	1.96	0.46
2:C:115:PHE:CE1	2:C:119:LEU:HD11	2.51	0.45
2:C:199:GLU:HG3	2:C:200:GLU:N	2.31	0.45
1:D:134:GLU:CD	1:D:134:GLU:H	2.19	0.45
1:D:5:ASP:O	1:D:9:LEU:HB2	2.15	0.45
2:B:257:LEU:HD11	7:B:505:HOH:O	2.16	0.45
1:A:97:LYS:N	1:A:97:LYS:HD2	2.31	0.45
1:A:167:GLU:CG	7:A:546:HOH:O	2.60	0.45
1:A:27:GLU:OE1	2:C:157:ARG:NH2	2.50	0.45
2:B:20:ILE:CD1	2:B:36:THR:HG21	2.47	0.45
2:B:253:LEU:HD12	2:B:256:SER:HG	1.81	0.45
2:B:1:MET:HG3	2:B:2:GLN:HG2	1.98	0.45
1:D:212:CYS:C	1:D:214:GLY:N	2.70	0.45
2:B:162:VAL:O	2:B:163:ASP:HB2	2.17	0.45
2:B:130:ASP:O	2:B:134:ARG:HG3	2.17	0.45
1:D:178:HIS:HE1	7:D:362:HOH:O	1.99	0.45
2:C:21:THR:HG22	7:C:766:HOH:O	2.17	0.45
1:D:135:ARG:O	1:D:139:VAL:HG23	2.17	0.44
2:B:137:ARG:HD3	2:B:192:ILE:HG21	1.98	0.44
1:D:222:ASP:HB3	1:D:249:PRO:CD	2.47	0.44
1:A:132:PRO:O	1:A:135:ARG:N	2.46	0.44
2:B:4:TYR:CE2	2:B:8:ILE:HD11	2.52	0.44
1:D:263:ARG:HH11	1:D:263:ARG:HB2	1.82	0.44
1:D:7:TYR:O	1:D:11:LYS:HG2	2.18	0.44
2:B:237:GLU:HG2	2:B:241:GLU:OE2	2.16	0.44
1:A:149:SER:C	1:A:154:ALA:HB2	2.37	0.44
1:A:132:PRO:HG2	1:A:135:ARG:HB2	1.99	0.44
2:C:25:PRO:HG2	2:C:28:VAL:CG2	2.48	0.44
1:D:135:ARG:CG	1:D:135:ARG:HH11	2.20	0.44
2:C:44:ALA:HB2	2:C:183:HIS:CD2	2.52	0.44
1:D:135:ARG:NH1	1:D:135:ARG:CG	2.81	0.44
2:C:86:ASP:OD2	1:D:106:GLU:CG	2.63	0.44
2:C:31:PRO:HG3	2:C:100:TYR:CE1	2.53	0.43
1:D:212:CYS:HB3	1:D:268:ALA:HB2	1.99	0.43
1:D:269:GLN:HE22	1:D:288:ASN:HD21	1.64	0.43
2:C:5:TRP:CZ2	2:C:257:LEU:HB2	2.53	0.43
1:D:205:LYS:NZ	7:D:365:HOH:O	2.50	0.43
2:B:211:GLN:CD	2:B:257:LEU:HD23	2.39	0.43
2:B:203:LYS:HE3	2:B:243:LEU:HA	2.00	0.43
1:A:33:GLU:HG2	1:A:104:PHE:CE2	2.53	0.43
1:A:44:LYS:NZ	1:A:44:LYS:HB2	2.33	0.43
1:A:213:ILE:HA	1:A:216:LEU:HB3	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:245:LYS:HD3	1:A:250:LYS:NZ	2.33	0.43
2:B:8:ILE:HG23	2:B:49:LEU:HD12	1.99	0.43
1:A:88:MET:CE	1:A:160:VAL:HG21	2.48	0.43
1:A:78:MET:HE3	1:A:117:LEU:HB2	1.99	0.43
2:B:93:LYS:N	2:B:94:PRO:CD	2.82	0.43
2:C:59:ARG:HG2	7:C:635:HOH:O	2.19	0.43
1:A:27:GLU:HG3	1:D:27:GLU:HB3	2.01	0.43
1:A:219:VAL:HG22	1:A:248:TYR:CE2	2.53	0.43
1:A:276:HIS:HA	1:A:277:PRO:HD3	1.91	0.43
1:D:222:ASP:O	1:D:226:VAL:HG12	2.18	0.43
1:A:215:LEU:HD12	1:A:215:LEU:HA	1.93	0.43
1:D:209:PHE:CZ	1:D:213:ILE:HD13	2.54	0.43
1:D:168:VAL:O	1:D:168:VAL:HG23	2.19	0.42
1:A:247:THR:HB	1:A:249:PRO:HD2	2.00	0.42
2:B:18:LYS:O	2:B:18:LYS:HG2	2.19	0.42
1:D:189:VAL:HG21	1:D:210:ALA:HB2	2.00	0.42
1:A:96:GLY:N	7:A:730:HOH:O	2.38	0.42
2:B:82:LEU:HD21	2:B:105:GLU:HB2	2.01	0.42
1:D:256:LYS:HA	1:D:259:GLU:OE1	2.20	0.42
1:A:272:LEU:HA	1:A:275:PHE:CE1	2.54	0.42
1:A:216:LEU:O	1:A:220:VAL:HG23	2.20	0.42
1:D:179:HIS:HA	1:D:183:ALA:HB2	2.02	0.42
1:D:214:GLY:C	1:D:216:LEU:H	2.21	0.42
2:B:137:ARG:CD	2:B:192:ILE:HD13	2.47	0.42
1:D:248:TYR:C	1:D:250:LYS:H	2.23	0.42
1:A:47:ARG:HE	1:A:184:LEU:HD23	1.85	0.42
2:C:88:SER:CB	1:D:107:SER:H	2.33	0.42
1:D:152:LEU:C	1:D:152:LEU:HD23	2.40	0.42
2:B:35:LEU:HD12	2:B:108:THR:HG21	2.02	0.42
2:C:148:PRO:HA	2:C:152:ILE:HB	2.01	0.42
2:C:28:VAL:O	2:C:31:PRO:HD2	2.20	0.42
1:A:156:GLN:O	1:A:156:GLN:HG3	2.20	0.42
1:A:254:VAL:O	1:A:255:GLU:C	2.58	0.42
1:D:226:VAL:HG13	1:D:226:VAL:O	2.18	0.41
1:D:45:ARG:HB3	1:D:48:PRO:HG2	2.01	0.41
2:B:2:GLN:H	2:B:2:GLN:HG2	1.67	0.41
2:C:182:MET:O	2:C:185:CYS:HB3	2.20	0.41
2:B:25:PRO:C	2:B:27:THR:H	2.24	0.41
1:A:141:GLY:O	1:A:144:ALA:HB3	2.20	0.41
2:B:169:ASP:CG	1:D:36:ARG:HE	2.23	0.41
1:D:135:ARG:HD3	1:D:194:ILE:O	2.21	0.41
1:D:168:VAL:CG2	1:D:251:LEU:HD21	2.45	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:140:LEU:HD13	2:B:139:ILE:HG21	2.01	0.41
1:D:189:VAL:CG1	1:D:206:LEU:HB3	2.51	0.41
1:D:20:GLU:HG2	1:D:40:LEU:HD13	2.01	0.41
2:B:238:LEU:O	2:B:242:GLU:HG2	2.21	0.41
2:B:84:LEU:HD12	2:B:89:ARG:HB3	2.03	0.41
1:D:293:ARG:HG3	1:D:293:ARG:HH21	1.86	0.41
1:D:47:ARG:N	1:D:48:PRO:HD2	2.36	0.41
1:D:156:GLN:O	1:D:159:ASP:HB3	2.21	0.41
2:B:5:TRP:CZ2	2:B:253:LEU:HG	2.56	0.41
2:B:138:VAL:HG22	2:B:189:CYS:SG	2.61	0.41
2:B:181:GLU:HB2	7:B:508:HOH:O	2.21	0.41
1:D:155:GLY:O	1:D:159:ASP:HB2	2.20	0.41
2:C:138:VAL:O	2:C:142:ILE:HG13	2.21	0.41
2:C:92:SER:HB2	2:C:94:PRO:HD2	2.02	0.41
2:C:2:GLN:HG2	7:C:476:HOH:O	2.21	0.41
1:D:44:LYS:HB2	1:D:46:VAL:HG23	2.03	0.41
1:A:248:TYR:N	1:A:249:PRO:CD	2.84	0.40
2:C:139:ILE:HD13	1:D:140:LEU:HD13	2.03	0.40
2:C:148:PRO:HB3	1:D:115:ALA:HB1	2.04	0.40
2:C:76:ALA:HB2	2:C:151:MET:CE	2.52	0.40
1:D:4:PHE:CE1	1:D:5:ASP:OD1	2.75	0.40
2:C:140:ILE:HD13	1:D:126:ALA:HB2	2.03	0.40
1:D:75:ILE:O	1:D:78:MET:HG3	2.21	0.40
1:D:269:GLN:HE22	1:D:288:ASN:ND2	2.19	0.40
2:C:2:GLN:HB2	2:C:5:TRP:HD1	1.86	0.40
2:C:147:GLY:HA3	7:C:341:HOH:O	2.21	0.40
1:D:143:LEU:HD13	1:D:191:LEU:HD11	2.03	0.40
1:D:214:GLY:O	1:D:216:LEU:N	2.55	0.40
2:C:145:ALA:HB1	2:C:182:MET:HE3	2.03	0.40
2:C:93:LYS:HA	2:C:93:LYS:HE2	2.03	0.40
2:B:54:LEU:HD13	2:B:250:ASN:OD1	2.22	0.40
2:B:49:LEU:HD22	2:B:59:ARG:HG3	2.04	0.40
2:B:214:LEU:HD11	2:B:232:ILE:HG23	2.03	0.40
1:D:173:LEU:HD11	1:D:177:HIS:HE2	1.87	0.40
2:B:215:ARG:NH2	7:B:325:HOH:O	2.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	275/295 (93%)	252 (92%)	22 (8%)	1 (0%)	43	60
1	D	280/295 (95%)	252 (90%)	25 (9%)	3 (1%)	21	28
2	B	252/274 (92%)	218 (86%)	30 (12%)	4 (2%)	14	18
2	C	257/274 (94%)	224 (87%)	22 (9%)	11 (4%)	4	2
All	All	1064/1138 (94%)	946 (89%)	99 (9%)	19 (2%)	13	15

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	223	SER
2	B	227	ILE
2	C	94	PRO
2	C	196	ALA
2	C	198	GLU
2	C	238	LEU
1	D	165	MET
1	D	168	VAL
1	A	42	GLY
2	C	57	GLY
2	C	224	HIS
2	C	226	LEU
2	C	237	GLU
2	B	26	GLU
2	C	91	VAL
2	B	163	ASP
2	C	60	SER
2	C	163	ASP
1	D	286	LEU

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	222/234 (95%)	214 (96%)	8 (4%)	47	67
1	D	226/234 (97%)	223 (99%)	3 (1%)	80	92
2	B	199/214 (93%)	196 (98%)	3 (2%)	76	90
2	C	201/214 (94%)	194 (96%)	7 (4%)	48	68
All	All	848/896 (95%)	827 (98%)	21 (2%)	60	79

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	44	LYS
1	A	90	ASN
1	A	99	THR
1	A	228	LYS
1	A	245	LYS
1	A	272	LEU
1	A	278	HIS
2	B	131	ASP
2	B	163	ASP
2	B	211	GLN
2	C	21	THR
2	C	86	ASP
2	C	93	LYS
2	C	94	PRO
2	C	131	ASP
2	C	182	MET
2	C	250	ASN
1	D	63	SER
1	D	78	MET
1	D	134	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	ASN
1	A	179	HIS
1	A	186	GLN
1	A	211	ASN

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Mol	Chain	Res	Type
2	B	34	HIS
2	B	81	HIS
2	B	156	HIS
2	B	183	HIS
2	B	202	GLN
2	B	206	ASN
2	B	224	HIS
2	C	33	HIS
1	D	265	ASN
1	D	288	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 ⓘ

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

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	$\# Z > 2$
3	PPV	A	901	4	8,8,8	1.54	1 (12%)	13,13,13	2.04	1 (7%)
5	EDO	C	902	-	3,3,3	0.55	0	2,2,2	0.40	0
6	GPP	D	900	-	18,18,18	1.96	6 (33%)	25,25,25	1.30	3 (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
3	PPV	A	901	4	-	0/6/6/6	0/0/0/0
5	EDO	C	902	-	-	0/1/1/1	0/0/0/0
6	GPP	D	900	-	-	0/19/19/19	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	900	GPP	C6-C7	-3.85	1.39	1.50
6	D	900	GPP	C2-C3	3.19	1.39	1.32
6	D	900	GPP	PB-O1B	3.08	1.61	1.51
6	D	900	GPP	C1-C2	-2.79	1.39	1.49
6	D	900	GPP	PA-O1A	2.65	1.61	1.51
3	A	901	PPV	P1-OPP	2.40	1.64	1.60
6	D	900	GPP	C7-C8	2.28	1.39	1.32

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	PPV	P2-OPP-P1	-7.04	111.03	131.68
6	D	900	GPP	PA-O3A-PB	-3.72	120.76	131.68
6	D	900	GPP	C4-C3-C5	2.81	119.66	115.39
6	D	900	GPP	C10-C8-C9	2.04	119.81	114.62

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	279/295 (94%)	-0.29	3 (1%) 77 77	28, 42, 65, 95	0
1	D	284/295 (96%)	0.13	21 (7%) 14 12	30, 47, 109, 125	0
2	B	256/274 (93%)	0.20	21 (8%) 12 10	32, 50, 128, 145	0
2	C	259/274 (94%)	0.08	9 (3%) 42 39	30, 50, 97, 106	0
All	All	1078/1138 (94%)	0.03	54 (5%) 28 25	28, 46, 104, 145	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	250	ASN	8.6
2	C	127	ALA	7.4
1	D	251	LEU	7.3
2	B	248	GLY	6.6
2	B	259	ALA	6.3
1	D	258	LYS	5.9
2	B	226	LEU	5.8
1	D	257	SER	5.7
2	C	249	LYS	5.5
2	B	249	LYS	5.3
2	B	235	LEU	4.4
1	A	166	ALA	4.2
2	B	254	MET	3.8
2	B	251	ALA	3.6
2	B	231	ILE	3.6
2	C	126	PRO	3.4
1	D	255	GLU	3.4
1	D	292	TYR	3.4
2	C	89	ARG	3.4
1	D	165	MET	3.2
1	D	260	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	168	VAL	3.2
2	B	234	LYS	3.1
2	C	226	LEU	3.1
2	B	258	VAL	3.0
2	B	224	HIS	3.0
1	D	289	TYR	2.9
1	D	167	GLU	2.8
2	B	257	LEU	2.8
2	B	256	SER	2.8
1	D	230	SER	2.7
2	B	238	LEU	2.7
1	D	130	GLY	2.7
1	D	163	GLU	2.6
1	D	248	TYR	2.6
2	B	1	MET	2.5
1	D	290	ILE	2.5
1	D	252	ILE	2.5
1	D	250	LYS	2.5
1	A	165	MET	2.5
2	B	253	LEU	2.4
2	B	230	ASN	2.4
1	D	164	GLY	2.4
2	C	162	VAL	2.4
2	B	222	ASN	2.3
2	C	59	ARG	2.3
2	B	236	LYS	2.3
1	A	167	GLU	2.3
2	C	222	ASN	2.3
1	D	217	PHE	2.2
2	C	225	GLN	2.1
1	D	232	GLU	2.0
1	D	231	LYS	2.0
2	B	240	LEU	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(\AA^2)	Q<0.9
5	EDO	C	902	4/4	0.14	0.16	35,37,38,38	0
6	GPP	D	900	19/19	0.18	-0.01	53,61,68,69	0
4	MG	A	905	1/1	0.11	-0.36	61,61,61,61	0
3	PPV	A	901	9/9	0.11	-0.42	83,84,86,87	0
4	MG	A	906	1/1	0.09	-0.99	96,96,96,96	0
4	MG	D	904	1/1	0.09	-1.61	57,57,57,57	0
4	MG	D	903	1/1	0.06	-2.99	48,48,48,48	0

6.5 Other polymers

There are no such residues in this entry.