



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 04:45 PM GMT

PDB ID : 2E7I
Title : Crystal Structure of Sep-tRNA:Cys-tRNA Synthase from Archaeoglobus fulgidus
Authors : Fukunaga, R.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-01-10
Resolution : 3.00 Å(reported)

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

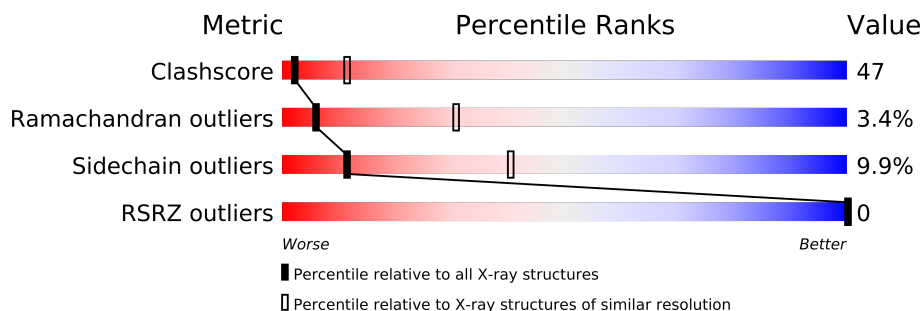
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.00 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	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

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 ≥ 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	371	
1	B	371	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 5586 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 Sep-tRNA:Cys-tRNA synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	0	0
			2737	1752	470	502	13			
1	B	344	Total	C	N	O	S	0	0	0
			2737	1752	470	502	13			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



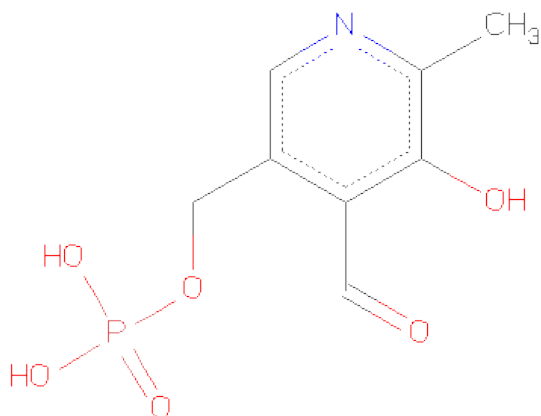
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 4 is water.

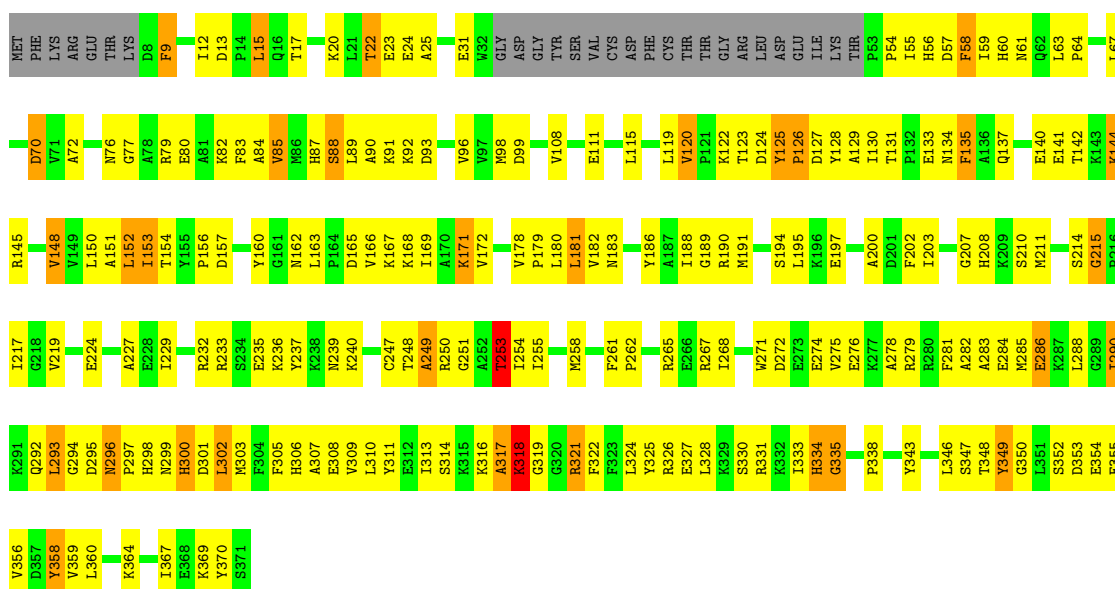
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	54	Total	O	0	0
			54	54		
4	B	18	Total	O	0	0
			18	18		

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.

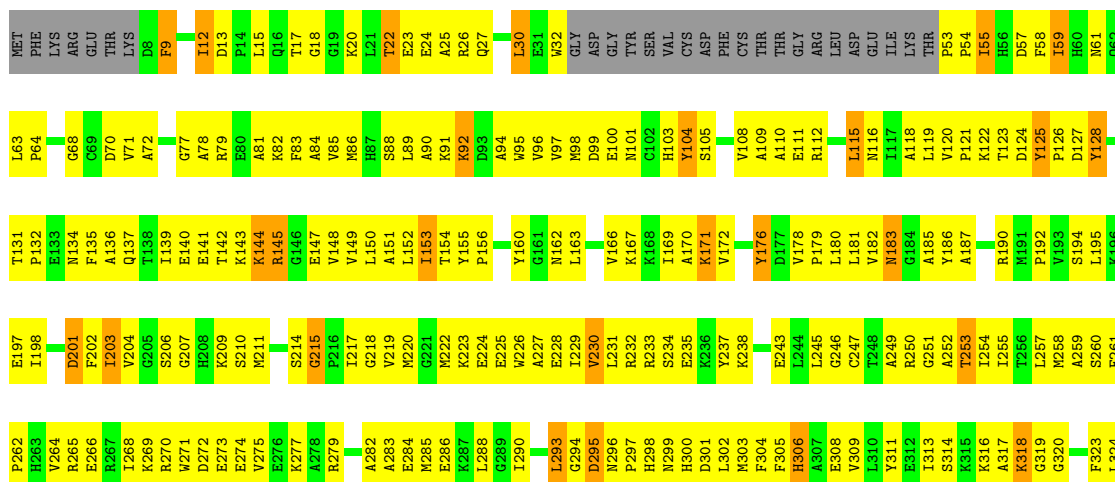
• Molecule 1: Sep-tRNA:Cys-tRNA synthase

Chain A:



• Molecule 1: Sep-tRNA:Cys-tRNA synthase

Chain B:





4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	220.66Å 220.66Å 92.93Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.44 – 3.00 47.44 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.44-3.00) 99.8 (47.44-3.00)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.271 , 0.318 0.283 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.383	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 31.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Outliers	0 of 27104 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5586	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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: SO4, PLP

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	0.39	0/2797	0.63	0/3769
1	B	0.33	0/2797	0.58	0/3769
All	All	0.36	0/5594	0.60	0/7538

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	2737	0	2745	246	0
1	B	2737	0	2747	276	0
2	A	15	0	0	0	0
2	B	10	0	0	0	0
3	A	15	0	6	5	0
4	A	54	0	0	8	0
4	B	18	0	0	3	0
All	All	5586	0	5498	521	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 47.

All (521) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:145:ARG:HH21	1:B:145:ARG:HB3	1.05	1.08
1:B:348:THR:HA	1:B:351:LEU:HD12	1.46	0.97
1:A:156:PRO:HB3	1:A:163:LEU:HD23	1.47	0.97
1:A:306:HIS:ND1	1:A:343:TYR:HE2	1.66	0.94
1:A:306:HIS:HD1	1:A:343:TYR:HE2	1.11	0.93
1:B:284:GLU:HG2	1:B:360:LEU:HD13	1.51	0.91
1:A:284:GLU:HG2	1:A:360:LEU:HD13	1.50	0.91
1:A:15:LEU:HD22	1:A:15:LEU:H	1.37	0.87
1:B:22:THR:HG22	1:B:25:ALA:H	1.35	0.87
1:A:22:THR:HG22	1:A:25:ALA:H	1.39	0.87
1:A:144:LYS:HE3	1:A:144:LYS:HA	1.56	0.86
1:B:145:ARG:NH2	1:B:145:ARG:HB3	1.89	0.86
1:B:97:VAL:HG12	1:B:118:ALA:HB3	1.55	0.85
1:B:318:LYS:HD2	1:B:319:GLY:H	1.41	0.85
1:B:181:LEU:HD21	1:B:204:VAL:HG23	1.60	0.83
1:A:64:PRO:HG2	4:A:1249:HOH:O	1.79	0.82
1:A:207:GLY:HA2	1:A:211:MET:HB2	1.60	0.81
1:B:83:PHE:CE2	1:B:112:ARG:HB3	2.15	0.81
1:A:325:TYR:CD1	1:A:338:PRO:HB3	2.16	0.81
1:A:137:GLN:O	1:A:141:GLU:HG3	1.80	0.80
1:A:9:PHE:CD2	1:A:334:HIS:HB3	2.16	0.80
1:A:22:THR:CG2	1:A:25:ALA:H	1.93	0.80
1:B:260:SER:O	1:B:264:VAL:HG23	1.81	0.80
1:B:71:VAL:HG23	1:B:222:MET:HG3	1.64	0.80
1:A:300:HIS:HD2	1:A:302:LEU:HB3	1.47	0.79
1:B:141:GLU:O	1:B:145:ARG:HG3	1.81	0.79
1:A:12:ILE:HD11	1:A:346:LEU:HD22	1.65	0.79
1:B:58:PHE:CE2	1:B:258:MET:HG2	2.19	0.78
1:B:123:THR:HG22	1:B:124:ASP:H	1.48	0.78
1:A:153:ILE:CD1	1:A:182:VAL:HG22	2.14	0.78
1:A:306:HIS:ND1	1:A:343:TYR:CE2	2.47	0.77
1:A:293:LEU:H	1:A:293:LEU:HD22	1.49	0.77
1:B:145:ARG:HH21	1:B:145:ARG:CB	1.92	0.76
1:A:99:ASP:HA	1:A:120:VAL:HG13	1.66	0.76
1:A:318:LYS:CD	1:A:319:GLY:H	1.99	0.76
1:B:192:PRO:HD3	1:B:268:ILE:HB	1.68	0.75
1:A:217:ILE:HB	1:A:253:THR:HG22	1.66	0.75
1:B:293:LEU:H	1:B:293:LEU:HD22	1.51	0.74
1:B:207:GLY:HA2	1:B:211:MET:HB2	1.68	0.73
1:A:318:LYS:HG2	1:A:319:GLY:N	2.04	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:154:THR:HA	1:B:183:ASN:HB3	1.71	0.72
1:B:12:ILE:HG23	1:B:335:GLY:HA3	1.70	0.72
1:B:290:ILE:O	1:B:290:ILE:HG13	1.88	0.72
1:B:104:TYR:HD1	1:B:105:SER:N	1.88	0.71
1:A:17:THR:HB	1:A:349:TYR:HB3	1.72	0.71
1:B:271:TRP:O	1:B:274:GLU:HB2	1.90	0.71
1:B:135:PHE:O	1:B:139:ILE:HG13	1.90	0.71
1:B:139:ILE:O	1:B:143:LYS:HG3	1.90	0.70
1:A:115:LEU:HA	4:A:1240:HOH:O	1.91	0.70
1:B:318:LYS:HD2	1:B:319:GLY:N	2.06	0.70
1:B:150:LEU:HD23	1:B:151:ALA:N	2.07	0.70
1:A:305:PHE:HE1	1:A:346:LEU:HD11	1.57	0.69
1:B:22:THR:HB	1:B:260:SER:OG	1.92	0.69
1:A:300:HIS:CD2	1:A:302:LEU:HB3	2.26	0.69
1:A:239:ASN:ND2	1:A:240:LYS:HE2	2.07	0.69
1:B:90:ALA:HA	1:B:115:LEU:HD21	1.75	0.69
1:A:194:SER:HB3	1:A:197:GLU:CG	2.21	0.69
1:B:79:ARG:HA	1:B:82:LYS:HE2	1.74	0.69
1:B:300:HIS:HD2	1:B:302:LEU:HB3	1.58	0.69
1:A:23:GLU:H	1:A:23:GLU:CD	1.95	0.69
1:B:136:ALA:HA	1:B:139:ILE:HD12	1.75	0.68
1:A:156:PRO:HB3	1:A:163:LEU:CD2	2.22	0.68
1:B:136:ALA:HB2	1:B:172:VAL:HG13	1.72	0.68
1:A:284:GLU:HB3	1:A:360:LEU:HD22	1.76	0.68
1:B:314:SER:O	1:B:320:GLY:HA2	1.94	0.68
1:A:15:LEU:CD2	1:A:15:LEU:H	2.06	0.68
1:B:59:ILE:O	1:B:59:ILE:HD12	1.93	0.67
1:B:194:SER:HB3	1:B:197:GLU:HG2	1.76	0.67
1:B:266:GLU:HG3	1:B:269:LYS:HD2	1.76	0.67
1:B:123:THR:HG22	1:B:124:ASP:N	2.08	0.67
1:B:86:MET:SD	1:B:110:ALA:HA	2.35	0.67
1:A:150:LEU:HD12	1:A:179:PRO:HB2	1.77	0.66
1:A:318:LYS:CG	1:A:319:GLY:N	2.58	0.66
1:B:234:SER:HB3	1:B:237:TYR:O	1.95	0.66
1:B:149:VAL:O	1:B:179:PRO:HD2	1.96	0.66
1:B:61:ASN:O	1:B:64:PRO:HD2	1.95	0.65
1:B:318:LYS:HE2	1:B:370:TYR:OH	1.96	0.65
1:A:293:LEU:N	1:A:293:LEU:HD22	2.10	0.65
1:A:292:GLN:HB2	1:A:305:PHE:CE2	2.31	0.65
1:A:188:ILE:O	1:A:210:SER:HB3	1.96	0.65
1:B:325:TYR:CG	1:B:338:PRO:HB3	2.32	0.65
1:B:182:VAL:O	1:B:203:ILE:HD12	1.96	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:223:LYS:HB2	1:B:226:TRP:HD1	1.62	0.65
1:A:167:LYS:O	1:A:171:LYS:HE2	1.97	0.65
1:A:20:LYS:HZ2	1:A:350:GLY:HA3	1.61	0.65
1:A:355:GLU:O	1:A:359:VAL:HG23	1.97	0.65
1:A:135:PHE:HE2	1:A:153:ILE:HG23	1.61	0.64
1:B:17:THR:HG21	1:B:347:SER:HB3	1.78	0.64
1:A:279:ARG:HH21	1:A:297:PRO:HB2	1.63	0.64
1:B:247:CYS:C	1:B:249:ALA:H	2.00	0.64
1:A:265:ARG:HG3	1:A:265:ARG:HH21	1.63	0.64
1:A:15:LEU:HD22	1:A:15:LEU:N	2.08	0.64
1:B:283:ALA:O	1:B:286:GLU:HB2	1.98	0.64
1:B:94:ALA:HB3	1:B:115:LEU:HD22	1.80	0.64
1:A:194:SER:HB3	1:A:197:GLU:HG3	1.78	0.64
1:A:63:LEU:HB3	1:A:64:PRO:HD3	1.80	0.64
1:A:313:ILE:HG22	1:A:317:ALA:HB2	1.80	0.64
1:B:71:VAL:HG23	1:B:222:MET:CG	2.28	0.63
1:B:342:ARG:HG2	1:B:343:TYR:CD1	2.33	0.63
1:A:108:VAL:HA	1:A:111:GLU:OE1	1.99	0.63
1:B:226:TRP:O	1:B:230:VAL:HG23	1.99	0.63
1:A:236:LYS:HB3	1:A:237:TYR:CD1	2.34	0.63
1:B:104:TYR:CD1	1:B:105:SER:N	2.66	0.63
1:A:276:GLU:HB3	1:B:273:GLU:HG2	1.81	0.62
1:A:58:PHE:CD1	1:A:258:MET:HG2	2.33	0.62
1:A:235:GLU:OE1	1:A:235:GLU:HA	1.99	0.62
1:A:346:LEU:O	1:A:346:LEU:HD12	2.00	0.62
1:A:208:HIS:HE1	3:A:1209:PLP:O1P	1.81	0.62
1:A:180:LEU:HD23	1:A:200:ALA:HB2	1.81	0.62
1:B:94:ALA:HB3	1:B:115:LEU:CD2	2.30	0.62
1:B:78:ALA:O	1:B:81:ALA:N	2.32	0.62
1:A:99:ASP:HA	1:A:120:VAL:CG1	2.30	0.62
1:B:63:LEU:HD11	1:B:219:VAL:HG11	1.82	0.61
1:B:85:VAL:HA	1:B:88:SER:HB3	1.81	0.61
1:A:12:ILE:HD11	1:A:346:LEU:CD2	2.31	0.61
1:A:195:LEU:HD13	1:A:203:ILE:HB	1.81	0.61
1:B:136:ALA:HB2	1:B:172:VAL:CG1	2.30	0.61
1:B:160:TYR:HA	1:B:304:PHE:HB2	1.80	0.61
1:B:83:PHE:CD2	1:B:112:ARG:HB3	2.36	0.61
1:A:321:ARG:HG3	1:A:322:PHE:CD2	2.36	0.61
1:B:83:PHE:HB2	1:B:109:ALA:HB1	1.82	0.61
1:A:20:LYS:NZ	1:A:350:GLY:HA3	2.15	0.61
1:B:323:PHE:HE1	1:B:341:THR:HG21	1.66	0.61
1:B:22:THR:HG22	1:B:25:ALA:N	2.14	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:265:ARG:O	1:B:268:ILE:HG12	2.02	0.60
1:A:251:GLY:O	1:A:255:ILE:HG12	2.01	0.60
1:A:194:SER:O	1:A:197:GLU:HB2	2.01	0.60
1:B:163:LEU:HD12	1:B:299:ASN:HB2	1.82	0.60
1:A:84:ALA:O	1:A:88:SER:HB2	2.01	0.60
1:B:253:THR:CG2	1:B:254:ILE:N	2.64	0.60
1:A:22:THR:HG22	1:A:25:ALA:CB	2.32	0.60
1:B:167:LYS:HD3	1:B:167:LYS:O	2.02	0.60
1:B:9:PHE:CD2	1:B:334:HIS:HB3	2.37	0.60
1:A:318:LYS:CG	1:A:319:GLY:H	2.15	0.59
1:B:293:LEU:HD21	1:B:306:HIS:HB3	1.82	0.59
1:B:63:LEU:HD21	1:B:219:VAL:HG13	1.82	0.59
1:B:100:GLU:OE1	1:B:121:PRO:HA	2.02	0.59
1:B:194:SER:HB3	1:B:197:GLU:CG	2.31	0.59
1:B:95:TRP:N	1:B:147:GLU:O	2.35	0.59
1:A:9:PHE:HD2	1:A:334:HIS:HB3	1.62	0.59
1:A:300:HIS:CD2	1:A:302:LEU:H	2.21	0.59
1:A:123:THR:HG22	1:A:124:ASP:N	2.17	0.59
1:A:250:ARG:HH21	1:A:250:ARG:HG2	1.67	0.59
1:B:140:GLU:O	1:B:144:LYS:HD2	2.02	0.59
1:B:57:ASP:HA	1:B:61:ASN:HD22	1.68	0.59
1:A:295:ASP:HB2	1:A:298:HIS:ND1	2.18	0.59
1:B:89:LEU:HD21	1:B:226:TRP:CZ3	2.36	0.58
1:B:125:TYR:CE2	1:B:126:PRO:HB3	2.38	0.58
1:A:58:PHE:CD2	1:A:58:PHE:C	2.75	0.58
1:B:12:ILE:HD11	1:B:346:LEU:CD2	2.33	0.58
1:A:12:ILE:HD11	1:A:346:LEU:HB2	1.84	0.58
1:A:79:ARG:HH11	3:A:1209:PLP:H5A1	1.67	0.58
1:B:232:ARG:HH21	1:B:232:ARG:HG2	1.66	0.58
1:A:56:HIS:O	1:A:60:HIS:HB2	2.04	0.58
1:A:152:LEU:HD12	1:A:152:LEU:C	2.24	0.58
1:B:12:ILE:HD11	1:B:346:LEU:CB	2.34	0.58
1:B:98:MET:HE2	1:B:152:LEU:HD21	1.84	0.58
1:B:293:LEU:N	1:B:293:LEU:HD22	2.18	0.58
1:A:12:ILE:CG1	1:A:346:LEU:HB2	2.34	0.57
1:B:363:PHE:O	1:B:367:ILE:HG13	2.04	0.57
1:A:166:VAL:HG22	4:A:1226:HOH:O	2.03	0.57
1:A:15:LEU:HD13	1:A:186:TYR:OH	2.04	0.57
1:B:89:LEU:HD11	1:B:226:TRP:HZ3	1.69	0.57
1:B:293:LEU:HD21	1:B:306:HIS:CB	2.34	0.57
1:A:142:THR:O	1:A:145:ARG:N	2.36	0.57
1:B:86:MET:SD	1:B:110:ALA:CA	2.93	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:282:ALA:HB2	1:B:303:MET:SD	2.45	0.57
1:B:153:ILE:H	1:B:153:ILE:HD12	1.70	0.56
1:B:55:ILE:O	1:B:59:ILE:HG23	2.05	0.56
1:A:88:SER:C	1:A:89:LEU:HD12	2.26	0.56
1:B:27:GLN:HA	1:B:30:LEU:HG	1.87	0.56
1:B:23:GLU:H	1:B:23:GLU:CD	2.08	0.56
1:A:239:ASN:HD21	1:A:240:LYS:HE2	1.69	0.56
1:B:70:ASP:OD1	1:B:224:GLU:HB2	2.05	0.56
1:A:154:THR:HA	1:A:183:ASN:HB3	1.87	0.56
1:A:79:ARG:HA	1:A:82:LYS:HE3	1.88	0.56
1:B:88:SER:CB	1:B:230:VAL:HG13	2.36	0.56
1:B:247:CYS:C	1:B:249:ALA:N	2.59	0.56
1:A:123:THR:HG22	1:A:124:ASP:H	1.71	0.56
1:A:364:LYS:HB3	1:A:364:LYS:HZ2	1.70	0.56
1:B:151:ALA:O	1:B:180:LEU:HA	2.06	0.56
1:B:12:ILE:HG12	1:B:12:ILE:O	2.06	0.56
1:A:83:PHE:CZ	1:A:87:HIS:HB2	2.41	0.56
1:A:290:ILE:HB	1:A:307:ALA:HB2	1.88	0.55
1:B:125:TYR:CD2	1:B:126:PRO:HB3	2.42	0.55
1:B:12:ILE:HD11	1:B:346:LEU:HB2	1.87	0.55
1:A:300:HIS:HD2	1:A:302:LEU:H	1.55	0.55
1:B:22:THR:HG23	1:B:24:GLU:OE1	2.06	0.55
1:A:122:LYS:HE3	1:A:128:TYR:CD1	2.42	0.55
1:B:353:ASP:HB2	4:B:610:HOH:O	2.06	0.55
1:B:340:LEU:HD22	1:B:343:TYR:O	2.07	0.55
1:B:104:TYR:C	1:B:104:TYR:HD1	2.10	0.54
1:A:178:VAL:HG23	1:A:178:VAL:O	2.06	0.54
1:B:251:GLY:O	1:B:255:ILE:HG12	2.07	0.54
1:A:309:VAL:HG13	1:A:310:LEU:HD13	1.88	0.54
1:A:153:ILE:HD11	1:A:182:VAL:HG22	1.86	0.54
1:A:80:GLU:HA	1:A:80:GLU:OE1	2.07	0.54
1:B:104:TYR:C	1:B:104:TYR:CD1	2.80	0.54
1:A:76:ASN:O	1:A:80:GLU:HG3	2.08	0.54
1:B:295:ASP:HB3	1:B:298:HIS:HA	1.89	0.54
1:B:101:ASN:HA	4:B:607:HOH:O	2.08	0.54
1:B:99:ASP:HA	1:B:120:VAL:CG1	2.37	0.54
1:A:123:THR:CG2	1:A:127:ASP:HB2	2.38	0.54
1:A:153:ILE:HD13	1:A:182:VAL:HG22	1.90	0.54
1:A:217:ILE:HB	1:A:253:THR:CG2	2.35	0.54
1:B:15:LEU:H	1:B:15:LEU:HD22	1.72	0.54
1:A:144:LYS:CA	1:A:144:LYS:HE3	2.34	0.54
1:A:325:TYR:CG	1:A:338:PRO:HB3	2.43	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:83:PHE:HE2	1:B:112:ARG:HB3	1.71	0.53
1:B:156:PRO:HB3	1:B:163:LEU:CD2	2.38	0.53
1:A:254:ILE:HG23	1:A:255:ILE:N	2.23	0.53
1:A:299:ASN:O	1:A:300:HIS:HB2	2.07	0.53
1:B:82:LYS:HG2	1:B:204:VAL:HG21	1.91	0.53
1:A:96:VAL:HB	1:A:115:LEU:HD13	1.90	0.53
1:A:22:THR:HG22	1:A:25:ALA:N	2.17	0.53
1:B:293:LEU:CD2	1:B:293:LEU:H	2.20	0.53
1:A:285:MET:CE	1:A:290:ILE:HD13	2.38	0.53
1:A:186:TYR:HE1	3:A:1209:PLP:O3	1.92	0.53
1:B:12:ILE:CG1	1:B:346:LEU:HB2	2.38	0.53
1:A:283:ALA:HA	1:A:286:GLU:HB2	1.90	0.53
1:A:181:LEU:HA	1:A:202:PHE:HB2	1.90	0.53
1:B:86:MET:HE1	1:B:109:ALA:HB3	1.91	0.53
1:B:53:PRO:HB2	1:B:55:ILE:HG13	1.91	0.53
1:A:58:PHE:CE1	1:A:258:MET:HG2	2.43	0.53
1:B:71:VAL:HG21	1:B:227:ALA:HB2	1.89	0.53
1:B:27:GLN:HA	1:B:30:LEU:CD1	2.38	0.53
1:A:123:THR:HB	1:A:129:ALA:H	1.74	0.53
1:A:288:LEU:HD22	1:A:367:ILE:CD1	2.39	0.53
1:B:272:ASP:HA	1:B:275:VAL:HG12	1.91	0.53
1:B:96:VAL:HA	1:B:150:LEU:O	2.09	0.52
1:A:262:PRO:O	1:A:265:ARG:HB2	2.09	0.52
1:A:364:LYS:NZ	1:A:364:LYS:HB3	2.24	0.52
1:B:144:LYS:HE3	1:B:144:LYS:N	2.25	0.52
1:B:119:LEU:N	1:B:119:LEU:HD22	2.23	0.52
1:B:26:ARG:O	1:B:30:LEU:HG	2.10	0.52
1:A:191:MET:HB3	1:A:271:TRP:CE2	2.44	0.52
1:B:253:THR:HG23	1:B:254:ILE:N	2.23	0.52
1:B:72:ALA:HA	1:B:220:MET:O	2.09	0.52
1:B:98:MET:HA	1:B:152:LEU:O	2.10	0.52
1:B:61:ASN:C	1:B:64:PRO:HD2	2.30	0.52
1:A:166:VAL:HA	1:A:169:ILE:HB	1.91	0.52
1:A:190:ARG:HB3	1:A:271:TRP:CZ3	2.45	0.52
1:B:68:GLY:O	1:B:195:LEU:HB3	2.10	0.52
1:A:79:ARG:HG2	3:A:1209:PLP:O3P	2.10	0.52
1:A:67:LEU:HD22	1:A:188:ILE:HD13	1.90	0.52
1:B:285:MET:HB3	1:B:290:ILE:CG1	2.40	0.52
1:A:22:THR:HG22	1:A:25:ALA:HB3	1.92	0.51
1:B:227:ALA:O	1:B:231:LEU:HB2	2.09	0.51
1:A:190:ARG:O	1:A:268:ILE:HG22	2.09	0.51
1:B:201:ASP:HA	1:B:223:LYS:HE3	1.93	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:309:VAL:HG13	1:A:310:LEU:H	1.75	0.51
1:B:79:ARG:HA	1:B:82:LYS:CE	2.40	0.51
1:B:156:PRO:HG3	1:B:300:HIS:CE1	2.45	0.51
1:B:126:PRO:HB3	1:B:293:LEU:HD12	1.92	0.51
1:A:125:TYR:CE2	1:A:126:PRO:HB3	2.46	0.51
1:B:166:VAL:HG23	1:B:167:LYS:N	2.26	0.51
1:A:54:PRO:HB2	1:A:57:ASP:HB3	1.93	0.51
1:B:306:HIS:CE1	1:B:308:GLU:HB2	2.45	0.51
1:B:285:MET:CE	1:B:290:ILE:HD11	2.40	0.51
1:A:57:ASP:O	1:A:61:ASN:HB2	2.11	0.51
1:B:156:PRO:HD3	1:B:187:ALA:HB2	1.92	0.50
1:A:285:MET:HB3	1:A:290:ILE:HG12	1.93	0.50
1:A:288:LEU:HD22	1:A:367:ILE:HD12	1.94	0.50
1:A:296:ASN:HA	1:A:297:PRO:C	2.32	0.50
1:B:194:SER:O	1:B:198:ILE:HG13	2.11	0.50
1:A:358:TYR:C	1:A:358:TYR:CD1	2.84	0.50
1:B:194:SER:CB	1:B:197:GLU:HG2	2.40	0.50
1:A:236:LYS:HB3	1:A:237:TYR:CE1	2.47	0.50
1:A:237:TYR:N	1:A:237:TYR:CD1	2.80	0.50
1:A:250:ARG:HG2	1:A:250:ARG:NH2	2.26	0.50
1:A:316:LYS:HG3	1:A:317:ALA:N	2.25	0.50
1:B:327:GLU:HB3	1:B:366:ILE:HD11	1.94	0.50
1:B:348:THR:CA	1:B:351:LEU:HD12	2.29	0.50
1:B:190:ARG:HD3	1:B:301:ASP:OD1	2.11	0.50
1:A:125:TYR:HB2	1:A:128:TYR:OH	2.11	0.50
1:B:217:ILE:CD1	1:B:257:LEU:HD22	2.42	0.50
1:A:333:ILE:HD11	1:A:358:TYR:CE1	2.46	0.50
1:A:70:ASP:OD2	1:A:70:ASP:N	2.43	0.50
1:A:91:LYS:O	1:A:92:LYS:C	2.50	0.50
1:A:12:ILE:HD11	1:A:346:LEU:CB	2.42	0.49
1:B:181:LEU:C	1:B:181:LEU:HD13	2.32	0.49
1:A:265:ARG:HG3	1:A:265:ARG:NH2	2.26	0.49
1:B:232:ARG:NH2	1:B:232:ARG:HG2	2.26	0.49
1:B:311:TYR:HB2	1:B:342:ARG:CB	2.42	0.49
1:B:12:ILE:HD11	1:B:346:LEU:HD22	1.94	0.49
1:A:352:SER:O	1:A:355:GLU:N	2.45	0.49
1:A:67:LEU:HD11	1:A:219:VAL:HG12	1.94	0.49
1:A:12:ILE:HG12	1:A:12:ILE:O	2.12	0.49
1:B:217:ILE:HD12	1:B:253:THR:HG23	1.94	0.49
1:B:156:PRO:CD	1:B:187:ALA:HB2	2.43	0.49
1:A:283:ALA:O	1:A:286:GLU:HB2	2.13	0.49
1:B:327:GLU:HB3	1:B:366:ILE:CD1	2.43	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:9:PHE:CE2	1:A:334:HIS:HB3	2.47	0.49
1:B:342:ARG:HG2	1:B:343:TYR:HD1	1.75	0.49
1:B:160:TYR:HB2	1:B:162:ASN:OD1	2.11	0.49
1:A:233:ARG:NH2	1:A:233:ARG:HG3	2.27	0.49
1:B:206:SER:OG	1:B:209:LYS:HG2	2.12	0.49
1:B:126:PRO:CB	1:B:293:LEU:HD12	2.42	0.49
1:A:285:MET:HE2	1:A:290:ILE:HD13	1.95	0.49
1:A:311:TYR:O	1:A:314:SER:OG	2.25	0.49
1:B:254:ILE:O	1:B:258:MET:HG3	2.12	0.48
1:B:318:LYS:CD	1:B:319:GLY:N	2.76	0.48
1:A:153:ILE:HD12	1:A:153:ILE:H	1.78	0.48
1:A:122:LYS:HE3	1:A:128:TYR:HD1	1.76	0.48
1:A:12:ILE:CD1	1:A:346:LEU:HB2	2.43	0.48
1:B:181:LEU:HA	1:B:202:PHE:HB2	1.95	0.48
1:A:166:VAL:N	4:A:1226:HOH:O	2.45	0.48
1:A:13:ASP:HB2	1:A:335:GLY:HA2	1.96	0.48
1:B:71:VAL:CG2	1:B:222:MET:HG3	2.41	0.48
1:B:311:TYR:HB2	1:B:342:ARG:HB2	1.96	0.48
1:B:81:ALA:O	1:B:84:ALA:HB3	2.14	0.48
1:A:272:ASP:O	1:A:275:VAL:HG12	2.14	0.48
1:A:124:ASP:O	1:A:125:TYR:C	2.51	0.48
1:B:155:TYR:O	1:B:155:TYR:CD1	2.67	0.48
1:A:296:ASN:HD22	1:A:296:ASN:C	2.16	0.48
1:B:311:TYR:CD2	1:B:342:ARG:HD2	2.49	0.48
1:B:166:VAL:HG23	1:B:198:ILE:HG23	1.96	0.48
1:A:135:PHE:CE2	1:A:153:ILE:HG23	2.46	0.48
1:B:325:TYR:CZ	1:B:329:LYS:HE2	2.49	0.48
1:A:55:ILE:HG12	1:A:258:MET:HE2	1.95	0.48
1:B:131:THR:OG1	1:B:134:ASN:ND2	2.47	0.48
1:B:295:ASP:CB	1:B:298:HIS:HA	2.42	0.48
1:A:247:CYS:SG	1:A:249:ALA:HB2	2.54	0.48
1:B:88:SER:OG	1:B:230:VAL:HA	2.13	0.48
1:A:302:LEU:HD12	1:A:302:LEU:C	2.34	0.48
1:B:85:VAL:HG23	1:B:86:MET:N	2.29	0.47
1:B:58:PHE:C	1:B:58:PHE:CD1	2.87	0.47
1:B:192:PRO:CD	1:B:268:ILE:HB	2.42	0.47
1:A:125:TYR:C	1:A:125:TYR:CD2	2.87	0.47
1:B:358:TYR:CD1	1:B:358:TYR:C	2.87	0.47
1:B:180:LEU:N	1:B:201:ASP:OD1	2.46	0.47
1:A:153:ILE:HG13	1:A:180:LEU:HD11	1.96	0.47
1:A:168:LYS:O	1:A:172:VAL:HG23	2.14	0.47
1:A:182:VAL:HB	1:A:203:ILE:HD12	1.97	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:259:ALA:O	1:B:262:PRO:HD2	2.14	0.47
1:B:294:GLY:O	1:B:295:ASP:C	2.52	0.47
1:B:254:ILE:HD12	1:B:257:LEU:HD23	1.96	0.47
1:B:100:GLU:HB2	1:B:120:VAL:O	2.14	0.47
1:B:325:TYR:CD1	1:B:338:PRO:HB3	2.49	0.47
1:B:279:ARG:CZ	1:B:297:PRO:HB2	2.45	0.47
1:A:64:PRO:HG3	1:A:72:ALA:HB3	1.96	0.47
1:B:288:LEU:HD22	1:B:367:ILE:HD12	1.96	0.47
1:A:92:LYS:HB3	4:A:1233:HOH:O	2.15	0.47
1:A:233:ARG:HG3	1:A:233:ARG:HH21	1.79	0.47
1:B:53:PRO:C	1:B:55:ILE:H	2.18	0.47
1:B:12:ILE:CG2	1:B:335:GLY:HA3	2.41	0.47
1:B:121:PRO:HG2	1:B:134:ASN:HB3	1.97	0.47
1:B:57:ASP:O	1:B:61:ASN:HB2	2.14	0.47
1:B:150:LEU:HG	1:B:179:PRO:HB2	1.96	0.47
1:A:85:VAL:O	1:A:89:LEU:HD13	2.15	0.47
1:B:55:ILE:O	1:B:58:PHE:HB3	2.15	0.46
1:B:285:MET:HE2	1:B:363:PHE:CE2	2.50	0.46
1:A:131:THR:OG1	1:A:134:ASN:ND2	2.48	0.46
1:A:306:HIS:CG	1:A:343:TYR:HE2	2.31	0.46
1:B:285:MET:HB3	1:B:290:ILE:HG12	1.98	0.46
1:B:18:GLY:HA3	1:B:190:ARG:HH11	1.81	0.46
1:A:90:ALA:O	1:A:91:LYS:HG3	2.15	0.46
1:B:77:GLY:HA2	1:B:217:ILE:C	2.36	0.46
1:B:277:LYS:HE2	1:B:277:LYS:HB3	1.80	0.46
1:B:123:THR:CG2	1:B:124:ASP:H	2.24	0.46
1:B:153:ILE:HD11	1:B:182:VAL:HG22	1.98	0.46
1:A:12:ILE:HG12	1:A:346:LEU:HB2	1.98	0.46
1:B:13:ASP:HB2	1:B:335:GLY:HA2	1.97	0.46
1:B:156:PRO:HB3	1:B:163:LEU:HD23	1.97	0.46
1:A:157:ASP:OD2	1:A:162:ASN:HB2	2.16	0.46
1:B:97:VAL:HG12	1:B:118:ALA:CB	2.37	0.46
1:B:309:VAL:O	1:B:313:ILE:HG13	2.16	0.46
1:B:311:TYR:HB2	1:B:342:ARG:HA	1.98	0.46
1:A:180:LEU:HD23	1:A:200:ALA:CB	2.45	0.46
1:B:122:LYS:HD2	1:B:128:TYR:HD1	1.79	0.46
1:A:284:GLU:HA	1:A:284:GLU:OE1	2.16	0.46
1:A:123:THR:HG21	1:A:127:ASP:HB2	1.97	0.46
1:B:342:ARG:HG2	1:B:343:TYR:CE1	2.51	0.46
1:A:285:MET:HE3	1:A:290:ILE:HG12	1.97	0.46
1:A:296:ASN:ND2	1:A:296:ASN:C	2.70	0.45
1:B:54:PRO:O	1:B:55:ILE:C	2.55	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:183:ASN:C	1:B:185:ALA:H	2.19	0.45
1:A:331:ARG:O	1:A:333:ILE:HG13	2.16	0.45
1:A:232:ARG:HH21	1:A:232:ARG:HG2	1.81	0.45
1:B:300:HIS:CD2	1:B:302:LEU:H	2.35	0.45
1:A:160:TYR:HB2	1:A:162:ASN:OD1	2.16	0.45
1:A:279:ARG:HE	1:A:297:PRO:CB	2.30	0.45
1:A:99:ASP:HB2	1:A:130:ILE:HG13	1.98	0.45
1:A:318:LYS:HE2	1:A:370:TYR:OH	2.15	0.45
1:B:148:VAL:HG13	1:B:148:VAL:O	2.16	0.45
1:B:27:GLN:O	1:B:30:LEU:HB2	2.16	0.45
1:B:305:PHE:CE1	1:B:346:LEU:HD11	2.52	0.45
1:A:254:ILE:HG13	1:A:258:MET:HG3	1.98	0.45
1:B:272:ASP:HA	1:B:275:VAL:CG1	2.46	0.45
1:A:279:ARG:HE	1:A:297:PRO:HB3	1.82	0.45
1:B:235:GLU:HA	1:B:235:GLU:OE1	2.17	0.45
1:A:306:HIS:HA	1:A:343:TYR:HD2	1.80	0.45
1:B:89:LEU:HD11	1:B:226:TRP:CZ3	2.51	0.45
1:A:296:ASN:HD22	1:A:297:PRO:N	2.15	0.45
1:A:318:LYS:HD2	1:A:319:GLY:H	1.79	0.45
1:B:227:ALA:C	1:B:229:ILE:H	2.20	0.45
1:A:308:GLU:O	1:A:311:TYR:HB3	2.16	0.44
1:A:274:GLU:HA	1:A:274:GLU:OE1	2.17	0.44
1:B:150:LEU:CG	1:B:179:PRO:HB2	2.47	0.44
1:A:370:TYR:CD1	1:A:370:TYR:N	2.86	0.44
1:A:150:LEU:CD1	1:A:179:PRO:HB2	2.45	0.44
1:B:15:LEU:N	1:B:15:LEU:HD22	2.32	0.44
1:A:150:LEU:HG	1:A:151:ALA:N	2.32	0.44
1:B:290:ILE:O	1:B:290:ILE:CG1	2.62	0.44
1:A:306:HIS:HA	1:A:343:TYR:CD2	2.52	0.44
1:A:67:LEU:HD11	1:A:219:VAL:CG1	2.48	0.44
1:A:182:VAL:HB	1:A:203:ILE:CD1	2.47	0.44
1:B:306:HIS:HB2	1:B:343:TYR:CE2	2.52	0.44
1:B:171:LYS:NZ	1:B:198:ILE:O	2.50	0.44
1:B:143:LYS:HE2	1:B:148:VAL:CG1	2.46	0.44
1:B:370:TYR:N	1:B:370:TYR:CD1	2.86	0.44
1:A:279:ARG:O	1:A:282:ALA:HB3	2.18	0.44
1:B:91:LYS:O	1:B:94:ALA:HB2	2.17	0.44
1:A:328:LEU:C	1:A:330:SER:N	2.69	0.44
1:A:140:GLU:O	1:A:144:LYS:HD2	2.17	0.44
1:A:60:HIS:C	4:A:1249:HOH:O	2.56	0.44
1:A:352:SER:O	1:A:353:ASP:C	2.57	0.44
1:A:254:ILE:HG23	1:A:255:ILE:H	1.81	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:214:SER:O	1:A:215:GLY:O	2.36	0.44
1:B:262:PRO:HA	1:B:265:ARG:NH2	2.32	0.43
1:B:217:ILE:HD13	1:B:257:LEU:HD22	1.99	0.43
1:B:311:TYR:HB2	1:B:342:ARG:CA	2.48	0.43
1:B:12:ILE:CD1	1:B:346:LEU:HB2	2.49	0.43
1:A:141:GLU:O	1:A:144:LYS:HB2	2.19	0.43
1:A:275:VAL:HG23	1:A:299:ASN:HA	2.00	0.43
1:B:92:LYS:HD2	1:B:92:LYS:HA	1.84	0.43
1:B:166:VAL:O	1:B:170:ALA:N	2.46	0.43
1:A:288:LEU:HD23	1:A:288:LEU:HA	1.81	0.43
1:A:152:LEU:HD12	1:A:153:ILE:N	2.34	0.43
1:B:125:TYR:CD2	1:B:125:TYR:C	2.92	0.43
1:A:189:GLY:O	1:A:267:ARG:HD3	2.19	0.43
1:A:302:LEU:HD12	1:A:303:MET:N	2.32	0.43
1:A:369:LYS:HD3	1:A:370:TYR:HE1	1.83	0.43
1:A:166:VAL:HG23	1:A:167:LYS:N	2.33	0.43
1:B:86:MET:SD	1:B:110:ALA:HB2	2.59	0.43
1:B:269:LYS:HA	4:B:619:HOH:O	2.18	0.43
1:B:123:THR:CG2	1:B:127:ASP:HB2	2.49	0.43
1:B:190:ARG:HG2	1:B:271:TRP:CZ3	2.52	0.43
1:B:354:GLU:HG3	1:B:355:GLU:N	2.34	0.43
1:B:58:PHE:CD2	1:B:258:MET:HG2	2.54	0.43
1:A:227:ALA:C	1:A:229:ILE:H	2.22	0.43
1:A:227:ALA:C	1:A:229:ILE:N	2.72	0.43
1:A:326:ARG:HA	1:A:326:ARG:HD3	1.91	0.43
1:B:142:THR:HA	1:B:145:ARG:HG3	2.00	0.43
1:A:148:VAL:HG22	1:A:148:VAL:O	2.18	0.43
1:B:169:ILE:HG22	1:B:169:ILE:O	2.19	0.43
1:A:293:LEU:H	1:A:293:LEU:CD2	2.26	0.42
1:A:353:ASP:O	1:A:356:VAL:HB	2.19	0.42
1:B:9:PHE:CE1	1:B:332:LYS:HA	2.54	0.42
1:A:63:LEU:HD21	1:A:219:VAL:HG13	2.01	0.42
1:B:151:ALA:O	1:B:180:LEU:HD12	2.20	0.42
1:A:54:PRO:O	1:A:55:ILE:C	2.57	0.42
1:A:58:PHE:HD2	1:A:58:PHE:C	2.19	0.42
1:B:225:GLU:HG3	1:B:226:TRP:CD1	2.55	0.42
1:B:340:LEU:HA	1:B:340:LEU:HD23	1.92	0.42
1:B:270:ARG:O	1:B:271:TRP:C	2.57	0.42
1:A:58:PHE:HD1	1:A:258:MET:SD	2.43	0.42
1:B:32:TRP:CZ3	1:B:252:ALA:HB2	2.54	0.42
1:A:9:PHE:CE2	1:A:334:HIS:CB	3.02	0.42
1:B:285:MET:O	1:B:286:GLU:C	2.56	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:143:LYS:C	1:B:144:LYS:HE3	2.39	0.42
1:B:245:LEU:O	1:B:247:CYS:N	2.53	0.42
1:A:326:ARG:HG2	4:A:1215:HOH:O	2.20	0.42
1:B:132:PRO:HA	1:B:169:ILE:HG12	2.01	0.42
1:B:313:ILE:HA	1:B:316:LYS:HE2	2.01	0.42
1:B:172:VAL:O	1:B:176:TYR:HD2	2.01	0.42
1:B:219:VAL:HG12	1:B:220:MET:N	2.34	0.42
1:B:323:PHE:CE1	1:B:341:THR:HG21	2.50	0.42
1:A:296:ASN:CA	1:A:297:PRO:C	2.88	0.42
1:B:103:HIS:H	1:B:154:THR:HG21	1.85	0.42
1:A:327:GLU:O	1:A:331:ARG:HD2	2.20	0.42
1:B:233:ARG:HB3	1:B:238:LYS:O	2.19	0.42
1:B:285:MET:HE1	1:B:363:PHE:CZ	2.55	0.42
1:B:301:ASP:OD2	1:B:347:SER:OG	2.36	0.42
1:A:76:ASN:HA	1:A:76:ASN:HD22	1.65	0.42
1:B:178:VAL:HG23	1:B:178:VAL:O	2.19	0.42
1:B:214:SER:O	1:B:215:GLY:O	2.37	0.42
1:B:77:GLY:H	1:B:218:GLY:N	2.18	0.42
1:B:303:MET:HB2	1:B:305:PHE:HE1	1.85	0.41
1:B:78:ALA:O	1:B:79:ARG:C	2.58	0.41
1:B:96:VAL:HG12	1:B:116:ASN:O	2.20	0.41
1:A:261:PHE:CE1	1:A:265:ARG:HD3	2.55	0.41
1:A:125:TYR:CD2	1:A:126:PRO:N	2.88	0.41
1:B:27:GLN:HA	1:B:30:LEU:CG	2.49	0.41
1:A:77:GLY:O	1:A:80:GLU:HB2	2.19	0.41
1:B:313:ILE:HG23	1:B:370:TYR:HB3	2.02	0.41
1:B:86:MET:CE	1:B:109:ALA:HB3	2.50	0.41
1:A:276:GLU:O	1:A:279:ARG:HB2	2.20	0.41
1:B:91:LYS:O	1:B:92:LYS:C	2.57	0.41
1:A:281:PHE:CD1	1:A:281:PHE:C	2.93	0.41
1:A:301:ASP:O	1:A:347:SER:HA	2.21	0.41
1:A:22:THR:HG23	1:A:24:GLU:N	2.35	0.41
1:A:165:ASP:HA	4:A:1226:HOH:O	2.21	0.41
1:B:99:ASP:HA	1:B:120:VAL:HG13	2.02	0.41
1:A:133:GLU:CD	1:A:133:GLU:N	2.73	0.41
1:B:88:SER:HB2	1:B:243:GLU:OE1	2.20	0.41
1:A:346:LEU:C	1:A:346:LEU:HD12	2.41	0.41
1:B:364:LYS:NZ	1:B:364:LYS:HB3	2.34	0.41
1:B:334:HIS:ND1	1:B:334:HIS:C	2.73	0.41
1:B:250:ARG:HG2	1:B:251:GLY:N	2.36	0.41
1:B:15:LEU:H	1:B:15:LEU:CD2	2.34	0.41
1:B:15:LEU:HD13	1:B:186:TYR:OH	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:137:GLN:O	1:B:141:GLU:HG3	2.21	0.41
1:A:186:TYR:CE1	3:A:1209:PLP:O3	2.73	0.41
1:B:57:ASP:HA	1:B:61:ASN:ND2	2.32	0.41
1:A:98:MET:O	1:A:119:LEU:HA	2.20	0.41
1:A:278:ALA:HB2	1:A:348:THR:OG1	2.20	0.41
1:B:190:ARG:HG3	1:B:349:TYR:CE1	2.56	0.40
1:B:301:ASP:O	1:B:347:SER:HA	2.21	0.40
1:B:152:LEU:O	1:B:152:LEU:HD12	2.21	0.40
1:A:55:ILE:HG12	1:A:258:MET:CE	2.51	0.40
1:A:294:GLY:O	1:A:295:ASP:C	2.60	0.40
1:A:154:THR:HG22	1:A:183:ASN:OD1	2.21	0.40
1:B:261:PHE:CE1	1:B:265:ARG:HD3	2.55	0.40
1:A:268:ILE:H	1:A:268:ILE:HG12	1.69	0.40
1:A:144:LYS:CE	1:A:144:LYS:HA	2.39	0.40
1:A:275:VAL:O	1:A:279:ARG:HB2	2.21	0.40
1:B:108:VAL:HA	1:B:111:GLU:HB2	2.03	0.40
1:A:296:ASN:HD22	1:A:297:PRO:CA	2.34	0.40
1:A:171:LYS:HE2	1:A:171:LYS:HB2	1.94	0.40
1:A:125:TYR:C	1:A:125:TYR:HD2	2.24	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	340/371 (92%)	283 (83%)	48 (14%)	9 (3%)	8	39
1	B	340/371 (92%)	279 (82%)	47 (14%)	14 (4%)	4	24
All	All	680/742 (92%)	562 (83%)	95 (14%)	23 (3%)	6	31

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	249	ALA

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Mol	Chain	Res	Type
1	A	317	ALA
1	B	12	ILE
1	B	92	LYS
1	B	115	LEU
1	B	317	ALA
1	A	215	GLY
1	A	321	ARG
1	A	349	TYR
1	B	30	LEU
1	B	55	ILE
1	B	215	GLY
1	B	246	GLY
1	B	295	ASP
1	A	253	THR
1	B	128	TYR
1	A	300	HIS
1	A	318	LYS
1	B	210	SER
1	B	228	GLU
1	A	335	GLY
1	B	203	ILE
1	B	335	GLY

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	288/312 (92%)	255 (88%)	33 (12%)	8	32
1	B	288/312 (92%)	264 (92%)	24 (8%)	16	52
All	All	576/624 (92%)	519 (90%)	57 (10%)	11	40

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	PHE
1	A	15	LEU
1	A	22	THR

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Mol	Chain	Res	Type
1	A	31	GLU
1	A	58	PHE
1	A	59	ILE
1	A	70	ASP
1	A	85	VAL
1	A	88	SER
1	A	93	ASP
1	A	120	VAL
1	A	125	TYR
1	A	126	PRO
1	A	135	PHE
1	A	144	LYS
1	A	148	VAL
1	A	152	LEU
1	A	153	ILE
1	A	171	LYS
1	A	181	LEU
1	A	224	GLU
1	A	248	THR
1	A	253	THR
1	A	286	GLU
1	A	290	ILE
1	A	293	LEU
1	A	296	ASN
1	A	302	LEU
1	A	318	LYS
1	A	324	LEU
1	A	334	HIS
1	A	354	GLU
1	A	358	TYR
1	B	9	PHE
1	B	20	LYS
1	B	22	THR
1	B	59	ILE
1	B	104	TYR
1	B	125	TYR
1	B	144	LYS
1	B	145	ARG
1	B	153	ILE
1	B	171	LYS
1	B	176	TYR
1	B	183	ASN

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Mol	Chain	Res	Type
1	B	201	ASP
1	B	230	VAL
1	B	253	THR
1	B	293	LEU
1	B	296	ASN
1	B	306	HIS
1	B	318	LYS
1	B	324	LEU
1	B	334	HIS
1	B	346	LEU
1	B	354	GLU
1	B	370	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	61	ASN
1	A	76	ASN
1	A	87	HIS
1	A	134	ASN
1	A	208	HIS
1	A	296	ASN
1	A	300	HIS
1	B	16	GLN
1	B	76	ASN
1	B	134	ASN
1	B	292	GLN
1	B	296	ASN
1	B	299	ASN
1	B	300	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	$\# Z > 2$
3	PLP	A	1209	1	14,15,16	1.05	0	20,22,23	1.01	0
2	SO4	A	602	-	4,4,4	0.31	0	6,6,6	0.13	0
2	SO4	A	604	-	4,4,4	0.31	0	6,6,6	0.07	0
2	SO4	A	605	-	4,4,4	0.25	0	6,6,6	0.09	0
2	SO4	B	601	-	4,4,4	0.37	0	6,6,6	0.13	0
2	SO4	B	603	-	4,4,4	0.29	0	6,6,6	0.10	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
3	PLP	A	1209	1	-	0/6/6/8	0/1/1/1
2	SO4	A	602	-	-	0/0/0/0	0/0/0/0
2	SO4	A	604	-	-	0/0/0/0	0/0/0/0
2	SO4	A	605	-	-	0/0/0/0	0/0/0/0
2	SO4	B	601	-	-	0/0/0/0	0/0/0/0
2	SO4	B	603	-	-	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, 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	344/371 (92%)	-0.37	0 100 100	4, 38, 76, 115	0
1	B	344/371 (92%)	-0.12	0 100 100	29, 81, 116, 138	0
All	All	688/742 (92%)	-0.24	0 100 100	4, 59, 109, 138	0

There are no RSRZ outliers to report.

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
3	PLP	A	1209	15/16	0.18	0.54	42,61,73,76	0
2	SO4	A	605	5/5	0.17	0.34	101,111,113,116	0
2	SO4	B	601	5/5	0.25	0.09	45,52,74,77	0
2	SO4	A	602	5/5	0.15	-0.45	61,63,70,73	0
2	SO4	B	603	5/5	0.16	-0.66	75,78,79,84	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	A	604	5/5	0.11	-1.93	111,111,113,118	0

6.5 Other polymers ⓘ

There are no such residues in this entry.