



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2025 – 03:01 AM EDT

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 X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

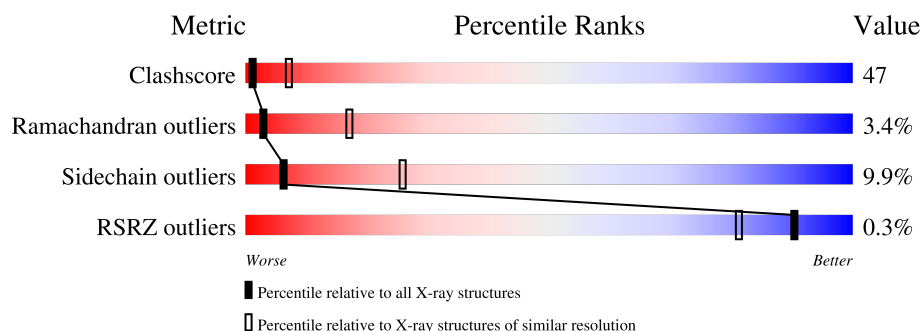
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

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	180529	2866 (3.00-3.00)
Ramachandran outliers	177936	2778 (3.00-3.00)
Sidechain outliers	177891	2781 (3.00-3.00)
RSRZ outliers	164620	2523 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

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 hydrogens and 0 are deuteriums.

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).



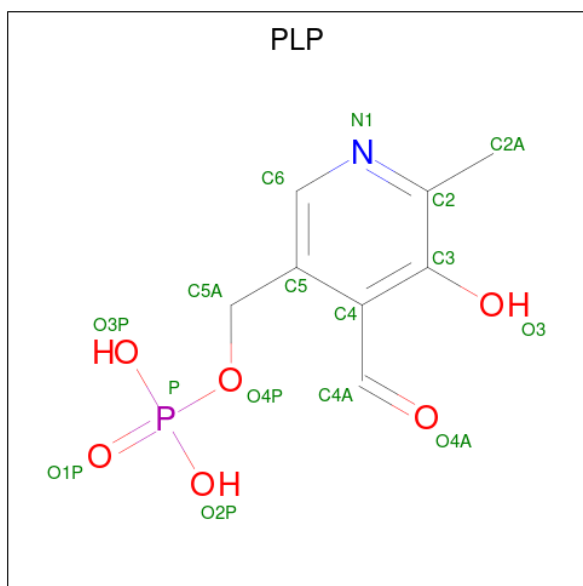
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $\text{C}_8\text{H}_{10}\text{NO}_6\text{P}$).



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

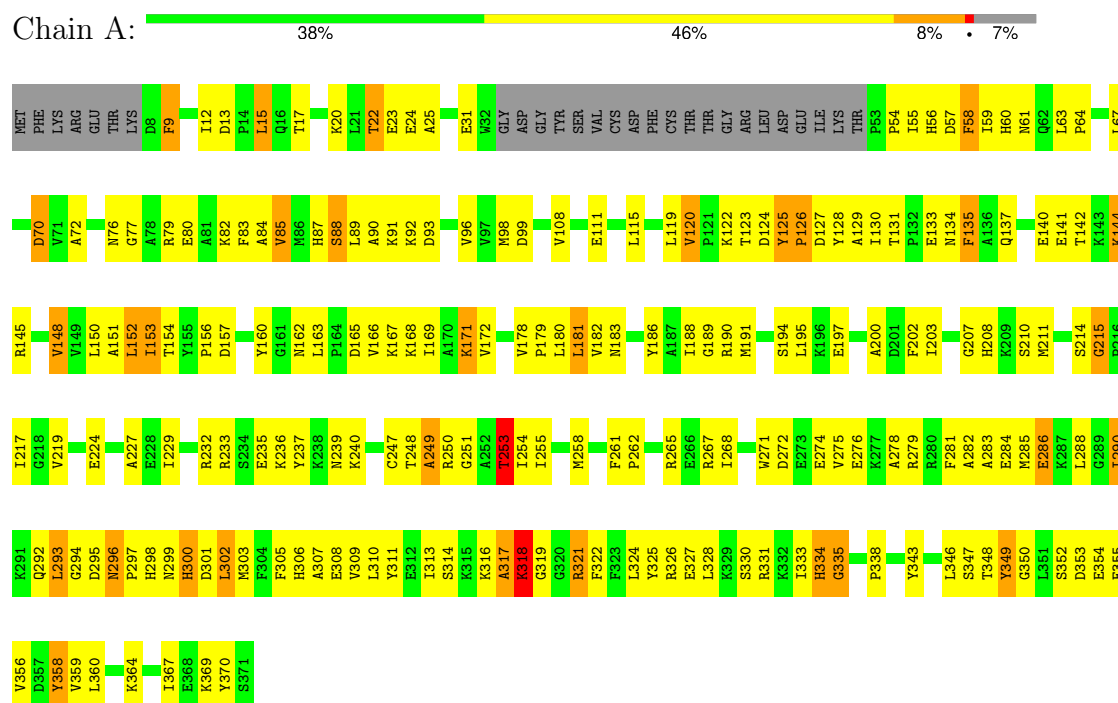
- Molecule 4 is water.

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

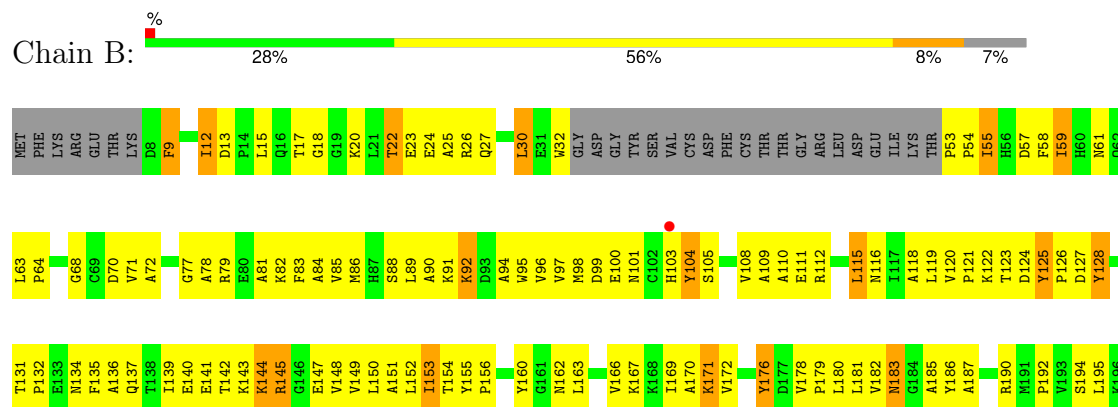
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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



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



E197	E198	D201	F202	I203	V204	G205	S206	G207	H208	K209	S210	M211	S214	G215	P216	I217	G218	V219	M220	G221	M222	K223	E224	E225	W226	A227	E228	I229	V230	L231	R232	R233	S234	E235	K236	Y237	K238	E243	L244	L245	G246	C247	T248	A249	R250	G251	A252	T253	I254	I255	T256	L257	M258	A259	S260	F261
P262	H263	V264	R265	E266	R267	I268	K269	R270	W271	D272	E273	E274	V275	E276	K277	A278	R279	A282	A283	E284	M285	E286	K287	L288	G289	I290	L293	G294	D295	N296	P297	H298	N299	H300	D301	L302	M303	F304	F305	H306	A307	E308	V309	L310	Y311	E312	I313	S314	K315	K316	A317	K318	G319	G320	F323	L324
Y325	R326	E327	L328	K329	K332	I333	H334	G335	P338	G339	L340	T341	R342	Y343	L346	S347	T348	Y349	G350	L351	S352	D353	E354	E355	Y358	Y359	L360	F363	K364	E365	I366	I367	Y370	S371																						

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 0.239 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.383	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 73.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	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.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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 Too-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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

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

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 47.

All (521) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (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:A:284:GLU:HG2	1:A:360:LEU:HD13	1.50	0.91
1:B:284:GLU:HG2	1:B:360:LEU:HD13	1.51	0.91
1:A:15:LEU:H	1:A:15:LEU:HD22	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:HB3	1:B:145:ARG:NH2	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:HD22	1:B:293:LEU:H	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	Interatomic distance (Å)	Clash overlap (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:HG13	1:B:290:ILE:O	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:A:239:ASN:ND2	1:A:240:LYS:HE2	2.07	0.69
1:A:300:HIS:CD2	1:A:302:LEU:HB3	2.26	0.69
1:B:22:THR:HB	1:B:260:SER:OG	1.92	0.69
1:A:194:SER:HB3	1:A:197:GLU:CG	2.21	0.69
1:B:90:ALA:HA	1:B:115:LEU:HD21	1.75	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:CD	1:A:23:GLU:H	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:H	1:A:15:LEU:CD2	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:A:293:LEU:HD22	1:A:293:LEU:N	2.10	0.65
1:B:318:LYS:HE2	1:B:370:TYR:OH	1.96	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:182:VAL:O	1:B:203:ILE:HD12	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325:TYR:CG	1:B:338:PRO:HB3	2.32	0.65
1:B:223:LYS:HB2	1:B:226:TRP:HD1	1.62	0.65
1:A:20:LYS:HZ2	1:A:350:GLY:HA3	1.61	0.65
1:A:167:LYS:O	1:A:171:LYS:HE2	1.97	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:265:ARG:HG3	1:A:265:ARG:HH21	1.63	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: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:A:194:SER:HB3	1:A:197:GLU:HG3	1.78	0.64
1:B:94:ALA:HB3	1:B:115:LEU:HD22	1.80	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:A:108:VAL:HA	1:A:111:GLU:OE1	1.99	0.63
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: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:HA	1:A:235:GLU:OE1	1.99	0.62
1:A:346:LEU:HD12	1:A:346:LEU:O	2.00	0.62
1:A:180:LEU:HD23	1:A:200:ALA:HB2	1.81	0.62
1:A:208:HIS:HE1	3:A:1209:PLP:O1P	1.81	0.62
1:B:94:ALA:HB3	1:B:115:LEU:CD2	2.30	0.62
1:A:99:ASP:HA	1:A:120:VAL:CG1	2.30	0.62
1:B:78:ALA:O	1:B:81:ALA:N	2.32	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:A:321:ARG:HG3	1:A:322:PHE:CD2	2.36	0.61
1:B:83:PHE:CD2	1:B:112:ARG:HB3	2.36	0.61
1:A:20:LYS:NZ	1:A:350:GLY:HA3	2.15	0.61
1:B:83:PHE:HB2	1:B:109:ALA:HB1	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:251:GLY:O	1:A:255:ILE:HG12	2.01	0.60
1:B:265:ARG:O	1:B:268:ILE:HG12	2.02	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:9:PHE:CD2	1:B:334:HIS:HB3	2.37	0.60
1:B:167:LYS:O	1:B:167:LYS:HD3	2.02	0.60
1:A:318:LYS:CG	1:A:319:GLY:H	2.15	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:293:LEU:HD21	1:B:306:HIS:HB3	1.82	0.59
1:B:95:TRP:N	1:B:147:GLU:O	2.35	0.59
1:B:194:SER:HB3	1:B:197:GLU:CG	2.31	0.59
1:A:9:PHE:HD2	1:A:334:HIS:HB3	1.62	0.59
1:A:123:THR:HG22	1:A:124:ASP:N	2.17	0.59
1:A:250:ARG:HG2	1:A:250:ARG:HH21	1.67	0.59
1:A:300:HIS:CD2	1:A:302:LEU:H	2.21	0.59
1:B:57:ASP:HA	1:B:61:ASN:HD22	1.68	0.59
1:B:140:GLU:O	1:B:144:LYS:HD2	2.02	0.59
1:A:295:ASP:HB2	1:A:298:HIS:ND1	2.18	0.59
1:A:58:PHE:C	1:A:58:PHE:CD2	2.75	0.58
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: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:HG2	1:B:232:ARG:HH21	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:HD22	1:B:293:LEU:N	2.18	0.58
1:A:12:ILE:CG1	1:A:346:LEU:HB2	2.34	0.57
1:A:166:VAL:HG22	4:A:1226:HOH:O	2.03	0.57
1:B:363:PHE:O	1:B:367:ILE:HG13	2.04	0.57
1:A:15:LEU:HD13	1:A:186:TYR:OH	2.04	0.57
1:A:142:THR:O	1:A:145:ARG:N	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:86:MET:SD	1:B:110:ALA:CA	2.93	0.57
1:B:282:ALA:HB2	1:B:303:MET:SD	2.45	0.57
1:B:153:ILE:HD12	1:B:153:ILE:H	1.70	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:55:ILE:O	1:B:59:ILE:HG23	2.05	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: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:247:CYS:C	1:B:249:ALA:N	2.59	0.56
1:A:83:PHE:CZ	1:A:87:HIS:HB2	2.41	0.56
1:B:12:ILE:HG12	1:B:12:ILE:O	2.06	0.56
1:B:151:ALA:O	1:B:180:LEU:HA	2.06	0.56
1:A:290:ILE:HB	1:A:307:ALA:HB2	1.88	0.55
1:B:12:ILE:HD11	1:B:346:LEU:HB2	1.87	0.55
1:B:125:TYR:CD2	1:B:126:PRO:HB3	2.42	0.55
1:A:300:HIS:HD2	1:A:302:LEU:H	1.55	0.55
1:A:122:LYS:HE3	1:A:128:TYR:CD1	2.42	0.55
1:B:22:THR:HG23	1:B:24:GLU:OE1	2.06	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:A:178:VAL:HG23	1:A:178:VAL:O	2.06	0.54
1:B:104:TYR:HD1	1:B:104:TYR:C	2.10	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:80:GLU:OE1	1:A:80:GLU:HA	2.07	0.54
1:A:153:ILE:HD11	1:A:182:VAL:HG22	1.86	0.54
1:A:76:ASN:O	1:A:80:GLU:HG3	2.08	0.54
1:B:104:TYR:CD1	1:B:104:TYR:C	2.80	0.54
1:B:101:ASN:HA	4:B:607:HOH:O	2.08	0.54
1:B:295:ASP:HB3	1:B:298:HIS:HA	1.89	0.54
1:A:123:THR:CG2	1:A:127:ASP:HB2	2.38	0.54
1:B:99:ASP:HA	1:B:120:VAL:CG1	2.37	0.54
1:A:153:ILE:HD13	1:A:182:VAL:HG22	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:ILE:HB	1:A:253:THR:CG2	2.35	0.54
1:B:15:LEU:HD22	1:B:15:LEU:H	1.72	0.54
1:A:144:LYS:HE3	1:A:144:LYS:CA	2.34	0.54
1:A:325:TYR:CG	1:A:338:PRO:HB3	2.43	0.54
1:A:254:ILE:HG23	1:A:255:ILE:N	2.23	0.53
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:299:ASN:O	1:A:300:HIS:HB2	2.07	0.53
1:A:96:VAL:HB	1:A:115:LEU:HD13	1.90	0.53
1:B:82:LYS:HG2	1:B:204:VAL:HG21	1.91	0.53
1:A:22:THR:HG22	1:A:25:ALA:N	2.17	0.53
1:A:285:MET:CE	1:A:290:ILE:HD13	2.38	0.53
1:B:293:LEU:H	1:B:293:LEU:CD2	2.20	0.53
1:A:186:TYR:HE1	3:A:1209:PLP:O3	1.92	0.53
1:A:283:ALA:HA	1:A:286:GLU:HB2	1.90	0.53
1:B:12:ILE:CG1	1:B:346:LEU:HB2	2.38	0.53
1:A:181:LEU:HA	1:A:202:PHE:HB2	1.90	0.53
1:A:58:PHE:CE1	1:A:258:MET:HG2	2.43	0.53
1:B:53:PRO:HB2	1:B:55:ILE:HG13	1.91	0.53
1:B:86:MET:HE1	1:B:109:ALA:HB3	1.91	0.53
1:B:27:GLN:HA	1:B:30:LEU:CD1	2.38	0.53
1:B:71:VAL:HG21	1:B:227:ALA:HB2	1.89	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:A:262:PRO:O	1:A:265:ARG:HB2	2.09	0.52
1:B:96:VAL:HA	1:B:150:LEU:O	2.09	0.52
1:A:364:LYS:HB3	1:A:364:LYS:NZ	2.24	0.52
1:B:119:LEU:N	1:B:119:LEU:HD22	2.23	0.52
1:B:144:LYS:HE3	1:B:144:LYS:N	2.25	0.52
1:A:191:MET:HB3	1:A:271:TRP:CE2	2.44	0.52
1:B:26:ARG:O	1:B:30:LEU:HG	2.10	0.52
1:B:72:ALA:HA	1:B:220:MET:O	2.09	0.52
1:B:253:THR:HG23	1:B:254:ILE:N	2.23	0.52
1:B:98:MET:HA	1:B:152:LEU:O	2.10	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:61:ASN:C	1:B:64:PRO:HD2	2.30	0.52
1:B:68:GLY:O	1:B:195:LEU:HB3	2.10	0.52
1:A:67:LEU:HD22	1:A:188:ILE:HD13	1.90	0.52
1:A:79:ARG:HG2	3:A:1209:PLP:O3P	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:190:ARG:O	1:A:268:ILE:HG22	2.09	0.51
1:B:227:ALA:O	1:B:231:LEU:HB2	2.09	0.51
1:B:201:ASP:HA	1:B:223:LYS:HE3	1.93	0.51
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:A:125:TYR:CE2	1:A:126:PRO:HB3	2.46	0.51
1:B:126:PRO:HB3	1:B:293:LEU:HD12	1.92	0.51
1:A:54:PRO:HB2	1:A:57:ASP:HB3	1.93	0.51
1:B:166:VAL:HG23	1:B:167:LYS:N	2.26	0.51
1:A:57:ASP:O	1:A:61:ASN:HB2	2.11	0.51
1:B:285:MET:CE	1:B:290:ILE:HD11	2.40	0.51
1:B:306:HIS:CE1	1:B:308:GLU:HB2	2.45	0.51
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:B:156:PRO:HD3	1:B:187:ALA:HB2	1.92	0.50
1:A:296:ASN:HA	1:A:297:PRO:C	2.32	0.50
1:A:358:TYR:CD1	1:A:358:TYR:C	2.84	0.50
1:B:194:SER:O	1:B:198:ILE:HG13	2.11	0.50
1:A:236:LYS:HB3	1:A:237:TYR:CE1	2.47	0.50
1:A:237:TYR:CD1	1:A:237:TYR:N	2.80	0.50
1:A:250:ARG:HG2	1:A:250:ARG:NH2	2.26	0.50
1:B:194:SER:CB	1:B:197:GLU:HG2	2.40	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:190:ARG:HD3	1:B:301:ASP:OD1	2.11	0.50
1:B:348:THR:CA	1:B:351:LEU:HD12	2.29	0.50
1:A:125:TYR:HB2	1:A:128:TYR:OH	2.11	0.50
1:A:70:ASP:N	1:A:70:ASP:OD2	2.43	0.50
1:A:91:LYS:O	1:A:92:LYS:C	2.50	0.50
1:A:333:ILE:HD11	1:A:358:TYR:CE1	2.46	0.50
1:B:217:ILE:CD1	1:B:257:LEU:HD22	2.42	0.50
1:A:12:ILE:HD11	1:A:346:LEU:CB	2.42	0.49
1:A:265:ARG:HG3	1:A:265:ARG:NH2	2.26	0.49
1:B:181:LEU:C	1:B:181:LEU:HD13	2.32	0.49
1:B:232:ARG:HG2	1:B:232:ARG:NH2	2.26	0.49
1:A:352:SER:O	1:A:355:GLU:N	2.45	0.49
1:B:12:ILE:HD11	1:B:346:LEU:HD22	1.94	0.49
1:B:311:TYR:HB2	1:B:342:ARG:CB	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:ILE:HG12	1:A:12:ILE:O	2.12	0.49
1:A:67:LEU:HD11	1:A:219:VAL:HG12	1.94	0.49
1:B:156:PRO:CD	1:B:187:ALA:HB2	2.43	0.49
1:B:217:ILE:HD12	1:B:253:THR:HG23	1.94	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
1:A:9:PHE:CE2	1:A:334:HIS:HB3	2.47	0.49
1:A:233:ARG:NH2	1:A:233:ARG:HG3	2.27	0.49
1:B:160:TYR:HB2	1:B:162:ASN:OD1	2.11	0.49
1:B:342:ARG:HG2	1:B:343:TYR:HD1	1.75	0.49
1:B:206:SER:OG	1:B:209:LYS:HG2	2.12	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:126:PRO:CB	1:B:293:LEU:HD12	2.42	0.49
1:B:254:ILE:O	1:B:258:MET:HG3	2.12	0.48
1:A:153:ILE:HD12	1:A:153:ILE:H	1.78	0.48
1:B:318:LYS:CD	1:B:319:GLY:N	2.76	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:A:13:ASP:HB2	1:A:335:GLY:HA2	1.96	0.48
1:A:166:VAL:N	4:A:1226:HOH:O	2.45	0.48
1:B:181:LEU:HA	1:B:202:PHE:HB2	1.95	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:A:124:ASP:O	1:A:125:TYR:C	2.51	0.48
1:A:272:ASP:O	1:A:275:VAL:HG12	2.14	0.48
1:B:81:ALA:O	1:B:84:ALA:HB3	2.14	0.48
1:B:155:TYR:O	1:B:155:TYR:CD1	2.67	0.48
1:A:296:ASN:C	1:A:296:ASN:HD22	2.16	0.48
1:B:166:VAL:HG23	1:B:198:ILE:HG23	1.96	0.48
1:B:311:TYR:CD2	1:B:342:ARG:HD2	2.49	0.48
1:A:55:ILE:HG12	1:A:258:MET:HE2	1.95	0.48
1:A:135:PHE:CE2	1:A:153:ILE:HG23	2.46	0.48
1:A:247:CYS:SG	1:A:249:ALA:HB2	2.54	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:B:325:TYR:CZ	1:B:329:LYS:HE2	2.49	0.48
1:A:302:LEU:C	1:A:302:LEU:HD12	2.34	0.48
1:B:88:SER:OG	1:B:230:VAL:HA	2.13	0.48
1:A:125:TYR:C	1:A:125:TYR:CD2	2.87	0.47
1:B:58:PHE:CD1	1:B:58:PHE:C	2.87	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:VAL:HG23	1:B:86:MET:N	2.29	0.47
1:B:192:PRO:CD	1:B:268:ILE:HB	2.42	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
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:100:GLU:HB2	1:B:120:VAL:O	2.14	0.47
1:B:254:ILE:HD12	1:B:257:LEU:HD23	1.96	0.47
1:B:279:ARG:CZ	1:B:297:PRO:HB2	2.45	0.47
1:B:325:TYR:CD1	1:B:338:PRO:HB3	2.49	0.47
1:A:64:PRO:HG3	1:A:72:ALA:HB3	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:288:LEU:HD22	1:B:367:ILE:HD12	1.96	0.47
1:B:12:ILE:CG2	1:B:335:GLY:HA3	2.41	0.47
1:B:53:PRO:C	1:B:55:ILE:H	2.18	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:A:85:VAL:O	1:A:89:LEU:HD13	2.15	0.47
1:B:150:LEU:HG	1:B:179:PRO:HB2	1.96	0.47
1:A:131:THR:OG1	1:A:134:ASN:ND2	2.48	0.46
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:90:ALA:O	1:A:91:LYS:HG3	2.15	0.46
1:A:306:HIS:CG	1:A:343:TYR:HE2	2.31	0.46
1:B:18:GLY:HA3	1:B:190:ARG:HH11	1.81	0.46
1:B:285:MET:HB3	1:B:290:ILE:HG12	1.98	0.46
1:B:77:GLY:HA2	1:B:217:ILE:C	2.36	0.46
1:B:277:LYS:HB3	1:B:277:LYS:HE2	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:A:157:ASP:OD2	1:A:162:ASN:HB2	2.16	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:123:THR:HG21	1:A:127:ASP:HB2	1.97	0.46
1:A:284:GLU:OE1	1:A:284:GLU:HA	2.16	0.46
1:A:285:MET:HE3	1:A:290:ILE:HG12	1.97	0.46
1:B:342:ARG:HG2	1:B:343:TYR:CE1	2.51	0.46
1:A:232:ARG:HG2	1:A:232:ARG:HH21	1.81	0.45
1:A:296:ASN:C	1:A:296:ASN:ND2	2.70	0.45
1:A:331:ARG:O	1:A:333:ILE:HG13	2.16	0.45
1:B:54:PRO:O	1:B:55:ILE:C	2.55	0.45
1:B:183:ASN:C	1:B:185:ALA:H	2.19	0.45
1:A:160:TYR:HB2	1:A:162:ASN:OD1	2.16	0.45
1:B:300:HIS:CD2	1:B:302:LEU:H	2.35	0.45
1:A:99:ASP:HB2	1:A:130:ILE:HG13	1.98	0.45
1:A:279:ARG:HE	1:A:297:PRO:CB	2.30	0.45
1:A:318:LYS:HE2	1:A:370:TYR:OH	2.15	0.45
1:B:27:GLN:O	1:B:30:LEU:HB2	2.16	0.45
1:B:148:VAL:HG13	1:B:148:VAL:O	2.16	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:B:305:PHE:CE1	1:B:346:LEU:HD11	2.52	0.45
1:A:279:ARG:HE	1:A:297:PRO:HB3	1.82	0.45
1:B:235:GLU:OE1	1:B:235:GLU:HA	2.17	0.45
1:A:306:HIS:HA	1:A:343:TYR:HD2	1.80	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:89:LEU:HD11	1:B:226:TRP:CZ3	2.51	0.45
1:B:227:ALA:C	1:B:229:ILE:H	2.20	0.45
1:A:274:GLU:HA	1:A:274:GLU:OE1	2.17	0.44
1:A:308:GLU:O	1:A:311:TYR:HB3	2.16	0.44
1:A:150:LEU:CD1	1:A:179:PRO:HB2	2.45	0.44
1:A:370:TYR:CD1	1:A:370:TYR:N	2.86	0.44
1:B:15:LEU:HD22	1:B:15:LEU:N	2.32	0.44
1:B:150:LEU:CG	1:B:179:PRO:HB2	2.47	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: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:A:306:HIS:HA	1:A:343:TYR:CD2	2.52	0.44
1:B:171:LYS:NZ	1:B:198:ILE:O	2.50	0.44
1:B:306:HIS:HB2	1:B:343:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:LYS:HE2	1:B:148:VAL:CG1	2.46	0.44
1:A:279:ARG:O	1:A:282:ALA:HB3	2.18	0.44
1:A:328:LEU:C	1:A:330:SER:N	2.69	0.44
1:B:91:LYS:O	1:B:94:ALA:HB2	2.17	0.44
1:B:370:TYR:N	1:B:370:TYR:CD1	2.86	0.44
1:A:60:HIS:C	4:A:1249:HOH:O	2.56	0.44
1:A:140:GLU:O	1:A:144:LYS:HD2	2.17	0.44
1:A:214:SER:O	1:A:215:GLY:O	2.36	0.44
1:A:254:ILE:HG23	1:A:255:ILE:H	1.81	0.44
1:A:352:SER:O	1:A:353:ASP:C	2.57	0.44
1:B:262:PRO:HA	1:B:265:ARG:NH2	2.32	0.43
1:B:12:ILE:CD1	1:B:346:LEU:HB2	2.49	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: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:A:288:LEU:HA	1:A:288:LEU:HD23	1.81	0.43
1:B:92:LYS:HA	1:B:92:LYS:HD2	1.84	0.43
1:B:166:VAL:O	1:B:170:ALA:N	2.46	0.43
1:A:152:LEU:HD12	1:A:153:ILE:N	2.34	0.43
1:A:189:GLY:O	1:A:267:ARG:HD3	2.19	0.43
1:B:125:TYR:CD2	1:B:125:TYR:C	2.92	0.43
1:A:166:VAL:HG23	1:A:167:LYS:N	2.33	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: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: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:HD3	1:A:326:ARG:HA	1.91	0.43
1:B:58:PHE:CD2	1:B:258:MET:HG2	2.54	0.43
1:A:148:VAL:O	1:A:148:VAL:HG22	2.18	0.43
1:B:142:THR:HA	1:B:145:ARG:HG3	2.00	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:54:PRO:O	1:A:55:ILE:C	2.57	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:PHE:C	1:A:58:PHE:HD2	2.19	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: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:B:225:GLU:HG3	1:B:226:TRP:CD1	2.55	0.42
1:B:270:ARG:O	1:B:271:TRP:C	2.57	0.42
1:B:340:LEU:HD23	1:B:340:LEU:HA	1.92	0.42
1:A:9:PHE:CE2	1:A:334:HIS:CB	3.02	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: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:B:285:MET:O	1:B:286:GLU:C	2.56	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:313:ILE:HA	1:B:316:LYS:HE2	2.01	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:A:327:GLU:O	1:A:331:ARG:HD2	2.20	0.42
1:B:103:HIS:H	1:B:154:THR:HG21	1.85	0.42
1:B:233:ARG:HB3	1:B:238:LYS:O	2.19	0.42
1:A:76:ASN:HD22	1:A:76:ASN:HA	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: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: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:A:77:GLY:O	1:A:80:GLU:HB2	2.19	0.41
1:A:125:TYR:CD2	1:A:126:PRO:N	2.88	0.41
1:A:261:PHE:CE1	1:A:265:ARG:HD3	2.55	0.41
1:B:27:GLN:HA	1:B:30:LEU:CG	2.49	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:B:313:ILE:HG23	1:B:370:TYR:HB3	2.02	0.41
1:A:276:GLU:O	1:A:279:ARG:HB2	2.20	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:B:86:MET:CE	1:B:109:ALA:HB3	2.50	0.41
1:B:91:LYS:O	1:B:92:LYS:C	2.57	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:THR:HG23	1:A:24:GLU:N	2.35	0.41
1:A:133:GLU:CD	1:A:133:GLU:N	2.73	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:346:LEU:HD12	1:A:346:LEU:C	2.41	0.41
1:B:88:SER:HB2	1:B:243:GLU:OE1	2.20	0.41
1:B:364:LYS:NZ	1:B:364:LYS:HB3	2.34	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
1:B:250:ARG:HG2	1:B:251:GLY:N	2.36	0.41
1:B:334:HIS:ND1	1:B:334:HIS:C	2.73	0.41
1:A:98:MET:O	1:A:119:LEU:HA	2.20	0.41
1:A:186:TYR:CE1	3:A:1209:PLP:O3	2.73	0.41
1:A:278:ALA:HB2	1:A:348:THR:OG1	2.20	0.41
1:B:57:ASP:HA	1:B:61:ASN:ND2	2.32	0.41
1:B:137:GLN:O	1:B:141:GLU:HG3	2.21	0.41
1:B:152:LEU:O	1:B:152:LEU:HD12	2.21	0.40
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:A:55:ILE:HG12	1:A:258:MET:CE	2.51	0.40
1:A:154:THR:HG22	1:A:183:ASN:OD1	2.21	0.40
1:A:294:GLY:O	1:A:295:ASP:C	2.60	0.40
1:A:268:ILE:H	1:A:268:ILE:HG12	1.69	0.40
1:B:261:PHE:CE1	1:B:265:ARG:HD3	2.55	0.40
1:A:144:LYS:HA	1:A:144:LYS:CE	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:125:TYR:C	1:A:125:TYR:HD2	2.24	0.40
1:A:171:LYS:HE2	1:A:171:LYS:HB2	1.94	0.40
1:A:296:ASN:HD22	1:A:297:PRO:CA	2.34	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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%)	4	23
1	B	340/371 (92%)	279 (82%)	47 (14%)	14 (4%)	2	13
All	All	680/742 (92%)	562 (83%)	95 (14%)	23 (3%)	3	17

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	249	ALA
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%)	4	20
1	B	288/312 (92%)	264 (92%)	24 (8%)	9	34
All	All	576/624 (92%)	519 (90%)	57 (10%)	6	26

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
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

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Mol	Chain	Res	Type
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
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

Sometimes 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 [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	603	-	4,4,4	0.38	0	6,6,6	0.10	0
3	PLP	A	1209	1	15,15,16	1.06	0	21,22,23	1.02	1 (4%)
2	SO4	A	605	-	4,4,4	0.39	0	6,6,6	0.08	0
2	SO4	B	601	-	4,4,4	0.42	0	6,6,6	0.11	0
2	SO4	A	604	-	4,4,4	0.39	0	6,6,6	0.08	0
2	SO4	A	602	-	4,4,4	0.37	0	6,6,6	0.14	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PLP	A	1209	1	-	2/6/6/8	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1209	PLP	O4P-P-O1P	2.24	112.51	106.44

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1209	PLP	C5A-O4P-P-O3P
3	A	1209	PLP	C5A-O4P-P-O1P

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1209	PLP	5	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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%)	-0.73	0	100 100	4, 38, 76, 115	0
1	B	344/371 (92%)	-0.14	2 (0%)	85 71	29, 81, 116, 138	0
All	All	688/742 (92%)	-0.44	2 (0%)	90 81	4, 59, 109, 138	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	103	HIS	2.5
1	B	203	ILE	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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. 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	RSCC	RSR	B-factors(Å ²)	Q < 0.9
2	SO4	A	605	5/5	0.80	0.14	101,111,113,116	0
2	SO4	A	604	5/5	0.81	0.13	111,111,113,118	0
2	SO4	B	601	5/5	0.93	0.08	45,52,74,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	B	603	5/5	0.94	0.12	75,78,79,84	0
2	SO4	A	602	5/5	0.95	0.07	61,63,70,73	0
3	PLP	A	1209	15/16	0.96	0.09	42,61,73,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.