



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:31 AM UTC

PDB ID : 2E7J / pdb_00002e7j
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 : 2.40 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

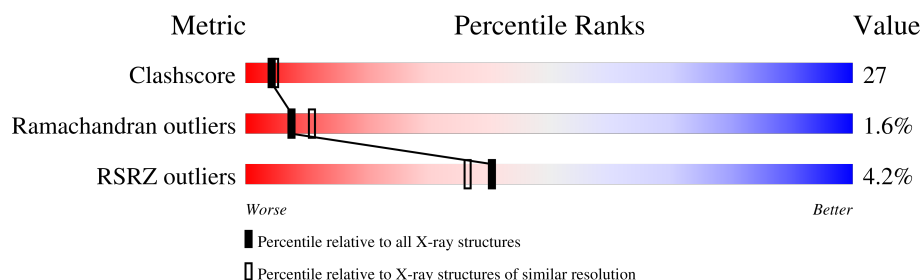
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 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	4% 44% 44% • 7%
1	B	371	4% 56% 33% • 7%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5605 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.

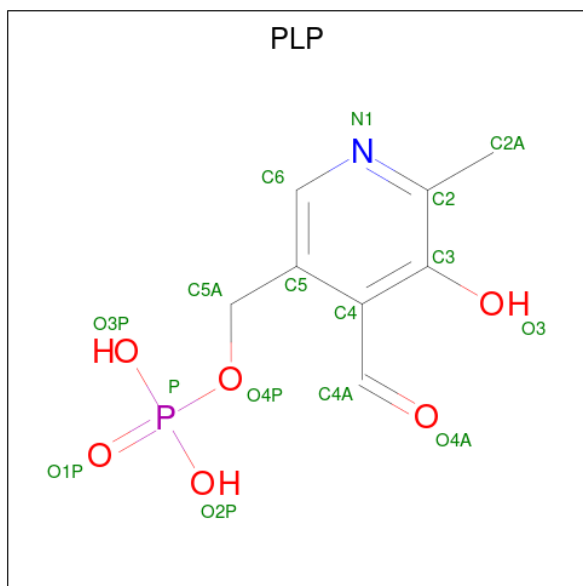
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	0	0
			2737	1752	470	502	13			
1	B	344	Total	C	N	O	S	0	0	0
			2737	1752	470	502	13			

- Molecule 2 is SULFATE ION (CCD ID: 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	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



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

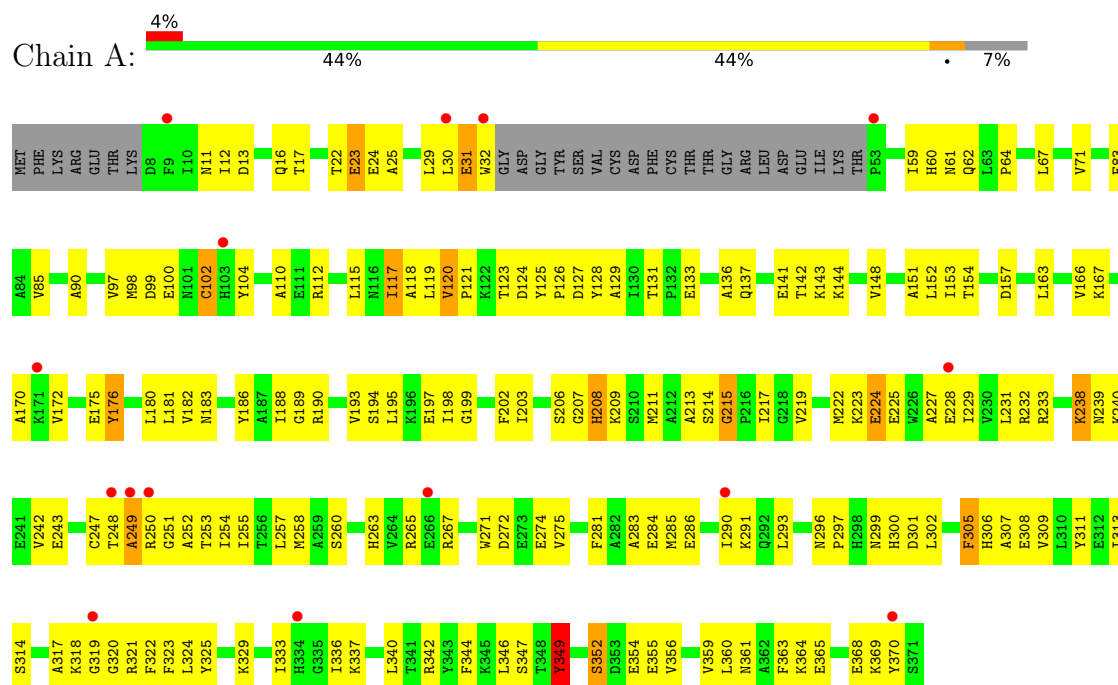
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	30	Total	O	0	0
			30	30		
4	B	51	Total	O	0	0
			51	51		

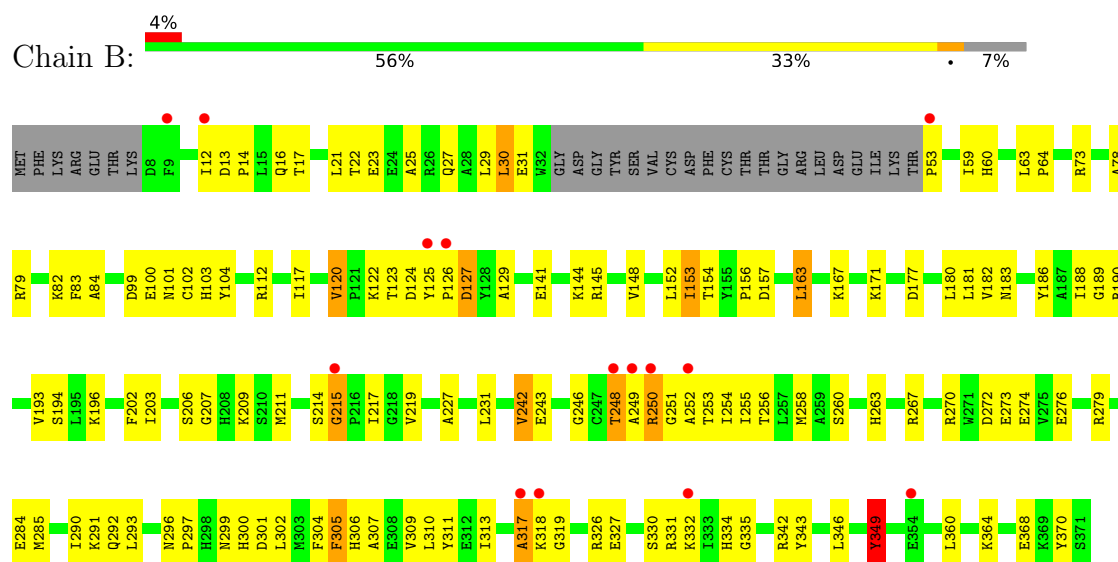
3 Residue-property plots

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	152.17Å 152.17Å 92.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.12 – 2.40 48.12 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.4 (48.12-2.40) 97.5 (48.12-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.202 , 0.255 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 59.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5605	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.69% 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	0.60	0/2797	1.07	15/3769 (0.4%)
1	B	0.67	0/2797	1.11	19/3769 (0.5%)
All	All	0.63	0/5594	1.09	34/7538 (0.5%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	THR	N-CA-C	-9.05	101.77	113.17
1	B	127	ASP	N-CA-C	8.30	119.96	111.07
1	A	117	ILE	N-CA-C	7.77	119.10	107.75
1	A	166	VAL	N-CA-C	7.31	117.44	110.42
1	B	120	VAL	N-CA-C	-7.05	102.57	108.63
1	A	291	LYS	N-CA-C	7.04	120.72	109.24
1	B	53	PRO	CA-C-N	6.74	128.26	119.84
1	B	53	PRO	C-N-CA	6.74	128.26	119.84
1	B	242	VAL	N-CA-C	6.71	116.86	110.42
1	A	23	GLU	N-CA-C	-6.59	104.02	111.14
1	B	291	LYS	N-CA-C	6.50	119.84	109.24
1	B	163	LEU	CA-C-N	-6.46	114.01	120.21
1	B	163	LEU	C-N-CA	-6.46	114.01	120.21
1	B	16	GLN	N-CA-C	6.39	122.15	113.97
1	B	273	GLU	N-CA-C	-6.12	104.52	111.07
1	B	305	PHE	N-CA-C	6.07	119.60	109.95
1	A	349	TYR	N-CA-C	6.05	123.69	110.80
1	B	177	ASP	CA-C-N	-6.00	118.11	123.33
1	B	177	ASP	C-N-CA	-6.00	118.11	123.33
1	A	62	GLN	N-CA-C	5.85	118.13	111.11
1	A	305	PHE	N-CA-C	5.84	118.99	109.59
1	B	153	ILE	CB-CA-C	-5.79	101.99	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	352	SER	N-CA-C	-5.64	102.93	110.55
1	B	317	ALA	N-CA-C	5.56	118.34	110.50
1	A	242	VAL	N-CA-C	5.52	116.16	110.36
1	B	248	THR	N-CA-C	-5.51	105.42	111.82
1	A	120	VAL	CA-C-N	5.49	125.79	119.92
1	A	120	VAL	C-N-CA	5.49	125.79	119.92
1	B	117	ILE	N-CA-C	5.26	115.54	108.17
1	A	102	CYS	N-CA-C	5.14	117.17	110.43
1	B	349	TYR	N-CA-C	5.07	121.60	110.80
1	B	124	ASP	N-CA-C	5.04	116.61	110.41
1	A	208	HIS	CA-C-N	-5.04	113.20	121.92
1	A	208	HIS	C-N-CA	-5.04	113.20	121.92

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	175	0
1	B	2737	0	2745	131	0
2	A	10	0	0	1	0
2	B	10	0	0	1	0
3	A	15	0	6	2	0
3	B	15	0	6	3	0
4	A	30	0	0	1	0
4	B	51	0	0	2	0
All	All	5605	0	5502	298	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 27.

All (298) 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:A:22:THR:HG22	1:A:25:ALA:H	1.03	1.05
1:A:102:CYS:HA	1:A:154:THR:HG21	1.41	0.99
1:B:217:ILE:HD12	1:B:253:THR:HG23	1.48	0.96
1:B:102:CYS:HA	1:B:154:THR:HG21	1.46	0.95
1:B:22:THR:HG22	1:B:25:ALA:H	1.33	0.92
1:B:318:LYS:HD2	1:B:319:GLY:N	1.86	0.90
1:A:318:LYS:HD2	1:A:319:GLY:N	1.85	0.89
1:A:22:THR:HG22	1:A:25:ALA:N	1.87	0.89
1:B:83:PHE:CE2	1:B:112:ARG:HG2	2.07	0.88
1:A:318:LYS:HD2	1:A:319:GLY:H	1.40	0.84
1:B:254:ILE:HG13	1:B:258:MET:HE2	1.61	0.82
1:B:12:ILE:HD11	1:B:346:LEU:HD22	1.63	0.80
1:B:327:GLU:O	1:B:331:ARG:HD2	1.83	0.79
1:B:217:ILE:HD12	1:B:253:THR:CG2	2.13	0.78
1:A:136:ALA:HB2	1:A:172:VAL:HG11	1.65	0.78
1:B:300:HIS:HD2	1:B:302:LEU:H	1.32	0.77
1:B:22:THR:CG2	1:B:25:ALA:H	1.98	0.76
1:A:300:HIS:HD2	1:A:302:LEU:H	1.33	0.76
1:A:337:LYS:HB3	1:A:340:LEU:HD12	1.68	0.76
1:A:300:HIS:CD2	1:A:302:LEU:H	2.05	0.74
1:A:23:GLU:CD	1:A:23:GLU:H	1.95	0.74
1:A:194:SER:HB3	1:A:197:GLU:HB2	1.70	0.74
1:A:285:MET:HB3	1:A:290:ILE:HG12	1.69	0.72
1:B:167:LYS:O	1:B:171:LYS:HE2	1.88	0.72
1:B:12:ILE:HD11	1:B:346:LEU:HD13	1.73	0.71
1:A:181:LEU:HD22	1:A:202:PHE:HB2	1.71	0.70
1:A:123:THR:HB	1:A:129:ALA:H	1.57	0.69
1:A:167:LYS:HE3	1:A:197:GLU:O	1.92	0.69
1:A:217:ILE:HD12	1:A:253:THR:HG23	1.73	0.69
1:B:12:ILE:CD1	1:B:346:LEU:HD22	2.22	0.69
1:A:233:ARG:HH21	1:A:233:ARG:HG3	1.58	0.69
1:A:284:GLU:HG2	1:A:360:LEU:HD13	1.73	0.69
1:B:100:GLU:OE2	1:B:122:LYS:HE2	1.92	0.69
1:A:83:PHE:CE2	1:A:112:ARG:HG2	2.29	0.68
1:A:136:ALA:HB2	1:A:172:VAL:CG1	2.23	0.68
1:B:22:THR:HG22	1:B:25:ALA:N	2.05	0.68
1:B:152:LEU:C	1:B:152:LEU:HD12	2.19	0.67
1:A:284:GLU:HB3	1:A:360:LEU:HD22	1.76	0.67
1:A:217:ILE:HD12	1:A:253:THR:CG2	2.24	0.67
1:B:194:SER:OG	1:B:196:LYS:HG2	1.93	0.67
1:B:318:LYS:HD2	1:B:319:GLY:H	1.60	0.67
1:A:228:GLU:C	1:A:229:ILE:HD13	2.20	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:GLU:OE2	1:B:112:ARG:HD2	1.96	0.66
1:B:251:GLY:O	1:B:255:ILE:HG12	1.96	0.65
1:B:300:HIS:CD2	1:B:302:LEU:H	2.13	0.65
1:A:22:THR:HB	1:A:260:SER:OG	1.95	0.65
1:A:254:ILE:HG13	1:A:258:MET:HE2	1.77	0.65
1:A:172:VAL:O	1:A:176:TYR:HD2	1.80	0.64
1:A:217:ILE:HB	1:A:253:THR:HG23	1.77	0.64
1:A:285:MET:HE3	1:A:290:ILE:CD1	2.26	0.64
1:B:250:ARG:HH21	1:B:250:ARG:HG2	1.61	0.64
1:B:27:GLN:O	1:B:30:LEU:HB2	1.97	0.63
1:A:12:ILE:HG12	1:A:12:ILE:O	1.98	0.63
1:B:250:ARG:HG2	1:B:250:ARG:NH2	2.14	0.63
1:A:285:MET:HE1	1:A:363:PHE:CD2	2.34	0.63
1:A:99:ASP:HA	1:A:120:VAL:HG13	1.81	0.62
1:A:271:TRP:O	1:A:275:VAL:HG12	1.99	0.62
1:A:318:LYS:HE3	1:A:323:PHE:HB2	1.81	0.62
1:A:100:GLU:OE1	1:A:121:PRO:HA	1.99	0.62
1:A:300:HIS:HD2	1:A:302:LEU:HB3	1.64	0.62
1:A:314:SER:O	1:A:320:GLY:HA2	2.00	0.61
1:A:300:HIS:CD2	1:A:302:LEU:HB3	2.36	0.61
1:B:30:LEU:O	1:B:31:GLU:C	2.43	0.61
1:B:148:VAL:HG13	1:B:148:VAL:O	2.01	0.61
1:A:285:MET:HE3	1:A:290:ILE:HD13	1.81	0.61
1:B:253:THR:HG22	1:B:254:ILE:N	2.16	0.60
1:A:194:SER:O	1:A:198:ILE:HG13	2.01	0.60
1:A:29:LEU:HD13	1:A:252:ALA:HB1	1.84	0.59
1:A:337:LYS:HB3	1:A:340:LEU:CD1	2.31	0.59
1:B:331:ARG:O	1:B:332:LYS:HB2	2.02	0.59
1:A:97:VAL:HG12	1:A:118:ALA:HB3	1.83	0.59
1:B:29:LEU:HD11	1:B:252:ALA:HB1	1.83	0.59
1:A:175:GLU:O	1:A:175:GLU:HG3	2.03	0.59
1:B:206:SER:OG	1:B:209:LYS:HG2	2.03	0.59
1:A:206:SER:OG	1:A:209:LYS:HG2	2.03	0.59
1:A:263:HIS:HE1	2:A:602:SO4:O3	1.86	0.59
1:B:189:GLY:O	1:B:267:ARG:HD3	2.03	0.58
1:A:67:LEU:HD11	1:A:219:VAL:CG1	2.33	0.58
1:B:17:THR:HB	1:B:349:TYR:HB3	1.84	0.58
1:B:22:THR:HG22	1:B:25:ALA:CB	2.33	0.58
1:B:13:ASP:HB2	1:B:335:GLY:HA2	1.86	0.57
1:A:112:ARG:HD2	1:B:243:GLU:OE2	2.04	0.57
1:A:11:ASN:HD21	1:A:16:GLN:HB2	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:THR:C	1:A:144:LYS:H	2.11	0.57
1:B:12:ILE:HG12	1:B:346:LEU:HB2	1.86	0.57
1:A:126:PRO:HB2	1:A:293:LEU:HD12	1.86	0.57
1:B:154:THR:OG1	1:B:157:ASP:HA	2.04	0.57
1:A:126:PRO:CB	1:A:293:LEU:HD12	2.35	0.56
1:A:17:THR:HB	1:A:349:TYR:HB3	1.86	0.56
1:B:22:THR:HB	1:B:260:SER:OG	2.06	0.56
1:A:12:ILE:CD1	1:A:346:LEU:HB2	2.36	0.56
1:A:99:ASP:OD1	1:A:154:THR:HG23	2.04	0.56
1:A:152:LEU:HD12	1:A:152:LEU:C	2.32	0.55
1:B:190:ARG:HD3	1:B:301:ASP:OD1	2.06	0.55
1:B:285:MET:HE3	1:B:290:ILE:CD1	2.36	0.55
1:A:61:ASN:O	1:A:64:PRO:HD2	2.06	0.55
1:B:346:LEU:HD12	1:B:346:LEU:C	2.32	0.55
1:A:232:ARG:HG2	1:A:232:ARG:HH21	1.71	0.55
1:B:29:LEU:CD1	1:B:252:ALA:HB1	2.37	0.55
1:A:352:SER:O	1:A:356:VAL:HG23	2.07	0.55
1:A:318:LYS:CD	1:A:319:GLY:N	2.67	0.54
1:B:12:ILE:HG12	1:B:12:ILE:O	2.06	0.54
1:B:194:SER:HG	1:B:196:LYS:HG2	1.71	0.54
1:B:188:ILE:CD1	1:B:193:VAL:HG21	2.38	0.54
1:B:217:ILE:CD1	1:B:253:THR:HG23	2.32	0.54
1:B:188:ILE:HD13	1:B:193:VAL:HG21	1.88	0.54
1:A:361:ASN:ND2	1:A:365:GLU:OE2	2.40	0.53
1:B:263:HIS:HE1	2:B:702:SO4:O2	1.92	0.53
1:A:142:THR:C	1:A:144:LYS:N	2.65	0.53
1:A:309:VAL:O	1:A:313:ILE:HG13	2.08	0.53
1:A:224:GLU:O	1:A:224:GLU:HG3	2.09	0.53
1:A:296:ASN:HD22	1:A:297:PRO:HA	1.74	0.53
1:A:32:TRP:CZ3	1:B:21:LEU:HD22	2.43	0.53
1:A:229:ILE:HD13	1:A:229:ILE:N	2.23	0.52
1:A:137:GLN:O	1:A:141:GLU:HG3	2.09	0.52
1:B:296:ASN:HD22	1:B:297:PRO:HA	1.73	0.52
1:A:213:ALA:HB2	1:A:257:LEU:HA	1.91	0.52
1:A:247:CYS:SG	1:A:249:ALA:HB2	2.49	0.52
1:A:175:GLU:O	1:A:175:GLU:CG	2.58	0.52
1:B:307:ALA:HB1	1:B:310:LEU:HD22	1.92	0.52
1:A:12:ILE:HG12	1:A:346:LEU:HB2	1.92	0.52
1:A:97:VAL:O	1:A:151:ALA:HA	2.09	0.51
1:B:152:LEU:HD12	1:B:153:ILE:N	2.25	0.51
1:B:156:PRO:HB3	1:B:163:LEU:HD23	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:PHE:HE1	1:A:346:LEU:HD11	1.76	0.51
1:B:326:ARG:O	1:B:330:SER:HB3	2.11	0.51
1:B:285:MET:HE3	1:B:290:ILE:HD13	1.92	0.51
1:A:265:ARG:HG3	1:A:265:ARG:HH21	1.75	0.51
1:B:84:ALA:HB1	1:B:243:GLU:HG3	1.93	0.51
1:B:123:THR:CG2	1:B:127:ASP:HB2	2.41	0.51
1:A:29:LEU:HD11	1:B:29:LEU:HD21	1.93	0.51
1:A:190:ARG:HD3	1:A:301:ASP:OD1	2.10	0.51
1:B:23:GLU:H	1:B:23:GLU:CD	2.19	0.51
1:A:29:LEU:HD12	1:A:29:LEU:O	2.11	0.50
1:A:30:LEU:O	1:A:31:GLU:C	2.53	0.50
1:B:104:TYR:CD1	1:B:104:TYR:C	2.89	0.50
1:A:181:LEU:HD13	1:A:182:VAL:N	2.25	0.50
1:A:285:MET:HE1	1:A:363:PHE:CE2	2.47	0.50
1:A:195:LEU:HD13	1:A:203:ILE:HB	1.92	0.50
1:A:143:LYS:HG3	1:A:148:VAL:HG11	1.92	0.50
1:A:274:GLU:OE1	1:A:349:TYR:O	2.30	0.50
1:A:251:GLY:O	1:A:255:ILE:HG12	2.12	0.49
1:B:279:ARG:HH21	1:B:299:ASN:ND2	2.10	0.49
1:A:188:ILE:HD11	1:A:203:ILE:HG13	1.94	0.49
1:A:250:ARG:HG2	1:A:251:GLY:N	2.27	0.49
1:A:207:GLY:HA2	1:A:211:MET:HB2	1.94	0.49
1:B:99:ASP:OD1	1:B:154:THR:HG23	2.12	0.49
1:B:59:ILE:HD13	1:B:73:ARG:HD3	1.93	0.49
1:B:186:TYR:O	1:B:300:HIS:HE1	1.95	0.49
1:B:250:ARG:HG2	1:B:251:GLY:H	1.78	0.49
1:B:99:ASP:HA	1:B:120:VAL:HG13	1.95	0.49
1:B:123:THR:HG22	1:B:127:ASP:HB2	1.93	0.49
1:B:318:LYS:CD	1:B:319:GLY:N	2.69	0.49
1:A:186:TYR:O	1:A:300:HIS:HE1	1.95	0.48
1:A:336:ILE:HD13	1:A:344:PHE:CD1	2.48	0.48
1:A:346:LEU:C	1:A:346:LEU:HD12	2.37	0.48
1:B:145:ARG:HG2	1:B:145:ARG:NH2	2.28	0.48
1:B:144:LYS:NZ	4:B:1244:HOH:O	2.46	0.48
1:B:207:GLY:HA2	1:B:211:MET:HB2	1.95	0.48
1:A:32:TRP:HZ3	1:B:21:LEU:HD22	1.79	0.48
1:A:112:ARG:HB2	1:B:242:VAL:CG1	2.44	0.48
1:A:239:ASN:HD21	1:A:240:LYS:HE2	1.78	0.48
1:A:222:MET:HE1	1:A:231:LEU:HD13	1.94	0.48
1:B:181:LEU:CD1	1:B:181:LEU:C	2.87	0.48
1:B:267:ARG:O	1:B:270:ARG:HG2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:ILE:HB	1:A:307:ALA:HB2	1.95	0.48
1:B:153:ILE:HD11	1:B:182:VAL:HG22	1.95	0.48
1:A:170:ALA:HB1	1:A:199:GLY:O	2.13	0.48
1:A:194:SER:HB3	1:A:197:GLU:CB	2.41	0.48
1:B:292:GLN:HB2	1:B:305:PHE:CE2	2.49	0.48
1:A:333:ILE:HD13	1:A:359:VAL:HG22	1.96	0.47
1:A:233:ARG:HD3	1:A:240:LYS:O	2.14	0.47
1:B:293:LEU:HD21	1:B:306:HIS:HB2	1.96	0.47
1:B:101:ASN:HB3	1:B:157:ASP:HB2	1.96	0.47
1:B:227:ALA:O	1:B:231:LEU:HB2	2.13	0.47
1:A:22:THR:CG2	1:A:24:GLU:HB2	2.44	0.47
1:A:153:ILE:HD12	1:A:153:ILE:H	1.79	0.47
1:A:214:SER:O	1:A:215:GLY:O	2.33	0.47
1:B:309:VAL:O	1:B:313:ILE:HG13	2.15	0.47
1:B:313:ILE:HD13	1:B:370:TYR:HB2	1.97	0.47
1:A:181:LEU:HD22	1:A:202:PHE:CB	2.41	0.47
1:A:153:ILE:HG13	1:A:180:LEU:HD11	1.96	0.47
1:A:285:MET:HE3	1:A:290:ILE:HD11	1.96	0.47
1:A:323:PHE:CG	1:A:324:LEU:N	2.82	0.47
1:B:214:SER:O	1:B:215:GLY:O	2.33	0.47
1:A:189:GLY:O	1:A:267:ARG:HD3	2.15	0.47
1:A:206:SER:HB3	1:A:209:LYS:CG	2.45	0.47
1:A:233:ARG:HG3	1:A:233:ARG:NH2	2.27	0.47
1:A:352:SER:OG	1:A:355:GLU:HG3	2.15	0.47
1:B:253:THR:CG2	1:B:254:ILE:N	2.78	0.47
1:A:233:ARG:NH2	1:A:233:ARG:CG	2.78	0.47
1:B:154:THR:HG1	1:B:157:ASP:HA	1.80	0.46
1:B:246:GLY:HA2	4:B:1224:HOH:O	2.15	0.46
1:B:153:ILE:HD12	1:B:153:ILE:H	1.81	0.46
1:A:12:ILE:HG13	1:A:347:SER:O	2.15	0.46
1:A:214:SER:C	1:A:215:GLY:O	2.55	0.46
1:B:59:ILE:HG13	1:B:60:HIS:CD2	2.50	0.46
1:A:12:ILE:HD11	1:A:346:LEU:HD22	1.97	0.46
1:B:141:GLU:O	1:B:144:LYS:HB2	2.16	0.46
1:B:252:ALA:O	1:B:256:THR:HG23	2.16	0.46
1:A:104:TYR:CD1	1:A:104:TYR:C	2.94	0.46
1:A:163:LEU:HD12	1:A:299:ASN:HB2	1.98	0.46
1:B:79:ARG:HA	1:B:82:LYS:HE2	1.97	0.46
1:A:13:ASP:O	1:A:17:THR:HG23	2.17	0.45
1:A:112:ARG:HB2	1:B:242:VAL:HG12	1.98	0.45
1:B:12:ILE:HD11	1:B:346:LEU:CD2	2.40	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:THR:HG21	1:A:24:GLU:HB2	1.97	0.45
1:B:318:LYS:HD2	1:B:319:GLY:CA	2.47	0.45
1:B:311:TYR:HB2	1:B:342:ARG:HG3	1.99	0.45
1:B:125:TYR:CD2	1:B:126:PRO:CA	3.00	0.45
1:B:214:SER:C	1:B:215:GLY:O	2.60	0.45
1:B:272:ASP:O	1:B:276:GLU:HG2	2.17	0.45
1:A:182:VAL:HB	1:A:203:ILE:HD12	1.98	0.45
1:B:206:SER:CB	1:B:209:LYS:HG2	2.47	0.45
1:A:206:SER:HB3	1:A:209:LYS:HG3	1.99	0.45
1:B:22:THR:HG22	1:B:25:ALA:HB2	1.99	0.45
1:B:274:GLU:OE1	1:B:349:TYR:O	2.34	0.45
1:A:131:THR:C	1:A:133:GLU:H	2.26	0.44
1:A:368:GLU:O	1:A:369:LYS:C	2.60	0.44
1:B:248:THR:O	1:B:249:ALA:HB3	2.17	0.44
1:A:12:ILE:CD1	1:A:346:LEU:HD22	2.48	0.44
1:A:209:LYS:NZ	3:A:1209:PLP:O3	2.50	0.44
1:B:78:ALA:HB3	3:B:1209:PLP:H5A2	1.98	0.44
1:A:99:ASP:CG	1:A:154:THR:HG23	2.43	0.44
1:A:153:ILE:HD11	1:A:182:VAL:HG22	2.00	0.44
1:A:313:ILE:O	1:A:317:ALA:HB2	2.18	0.44
1:B:125:TYR:HD2	1:B:126:PRO:N	2.15	0.44
1:B:183:ASN:C	1:B:183:ASN:ND2	2.76	0.44
1:A:123:THR:O	1:A:128:TYR:HA	2.16	0.44
1:B:181:LEU:HD22	1:B:202:PHE:HB2	1.99	0.43
1:B:313:ILE:HG22	1:B:317:ALA:HB2	1.98	0.43
1:B:12:ILE:HD11	1:B:346:LEU:CD1	2.47	0.43
1:B:12:ILE:HG22	1:B:334:HIS:O	2.18	0.43
1:A:123:THR:CG2	1:A:127:ASP:HB2	2.49	0.43
1:B:182:VAL:HB	1:B:203:ILE:HD12	2.00	0.43
1:B:305:PHE:O	1:B:343:TYR:HA	2.19	0.43
1:A:188:ILE:HG12	1:A:193:VAL:HG21	2.01	0.43
1:A:313:ILE:HD13	1:A:370:TYR:HB2	2.00	0.43
1:A:181:LEU:CD1	1:A:181:LEU:C	2.92	0.43
1:A:125:TYR:C	1:A:125:TYR:CD2	2.97	0.43
1:A:232:ARG:HG2	1:A:232:ARG:NH2	2.34	0.43
1:A:254:ILE:HG23	1:A:255:ILE:N	2.32	0.43
1:A:311:TYR:CE2	1:A:342:ARG:NH1	2.87	0.43
1:B:101:ASN:HB3	1:B:157:ASP:CB	2.49	0.42
1:B:250:ARG:HH21	1:B:250:ARG:CG	2.26	0.42
1:A:110:ALA:HB3	1:A:117:ILE:HD11	2.00	0.42
1:A:354:GLU:HG3	1:A:355:GLU:N	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:TYR:CD1	1:A:370:TYR:N	2.87	0.42
1:B:145:ARG:HG2	1:B:145:ARG:HH21	1.83	0.42
1:B:304:PHE:CZ	1:B:343:TYR:HB3	2.54	0.42
1:B:82:LYS:HZ1	3:B:1209:PLP:H6	1.83	0.42
1:A:321:ARG:HG3	1:A:322:PHE:CD2	2.55	0.42
1:B:63:LEU:HB3	1:B:64:PRO:HD3	2.02	0.42
1:A:120:VAL:O	1:A:120:VAL:HG22	2.18	0.42
1:A:154:THR:OG1	1:A:157:ASP:HA	2.18	0.42
1:A:154:THR:HA	1:A:183:ASN:HB3	2.00	0.42
1:A:22:THR:HG23	1:A:23:GLU:N	2.34	0.42
1:A:188:ILE:HD11	1:A:203:ILE:CG1	2.50	0.42
1:A:272:ASP:O	1:A:275:VAL:HG13	2.20	0.42
1:A:22:THR:HG23	1:A:24:GLU:H	1.85	0.42
1:A:305:PHE:CE1	1:A:346:LEU:HD11	2.55	0.42
1:B:284:GLU:HG2	1:B:360:LEU:HD13	2.01	0.42
1:A:324:LEU:HD23	1:A:324:LEU:HA	1.85	0.42
1:A:12:ILE:CG1	1:A:346:LEU:HB2	2.50	0.42
1:A:61:ASN:C	1:A:64:PRO:HD2	2.45	0.42
1:A:98:MET:O	1:A:119:LEU:HA	2.19	0.42
1:A:233:ARG:HD2	1:A:238:LYS:O	2.20	0.42
1:A:283:ALA:O	1:A:286:GLU:HB2	2.20	0.42
1:B:125:TYR:CD2	1:B:126:PRO:HA	2.55	0.42
1:B:145:ARG:HH21	1:B:145:ARG:CG	2.32	0.42
1:A:223:LYS:HB2	1:A:225:GLU:HG2	2.02	0.41
1:B:14:PRO:HB3	1:B:302:LEU:HD13	2.02	0.41
1:A:71:VAL:HG21	1:A:227:ALA:HB2	2.01	0.41
1:A:308:GLU:OE1	1:A:311:TYR:HD2	2.03	0.41
1:A:325:TYR:OH	1:A:329:LYS:HG3	2.21	0.41
1:A:250:ARG:HG2	1:A:250:ARG:HH21	1.85	0.41
1:A:206:SER:CB	1:A:209:LYS:HG2	2.50	0.41
1:A:293:LEU:HD21	1:A:306:HIS:HB2	2.02	0.41
1:B:183:ASN:C	1:B:183:ASN:HD22	2.29	0.41
1:B:364:LYS:O	1:B:368:GLU:HG3	2.20	0.41
1:A:281:PHE:CD2	1:A:360:LEU:HD21	2.56	0.41
1:A:124:ASP:OD1	1:A:125:TYR:N	2.52	0.41
1:A:364:LYS:O	1:A:368:GLU:HG3	2.20	0.41
1:B:153:ILE:HG13	1:B:180:LEU:HD11	2.03	0.41
1:A:85:VAL:HG21	1:A:181:LEU:HD23	2.03	0.41
1:A:12:ILE:HD11	1:A:346:LEU:HB2	2.03	0.41
1:A:32:TRP:CH2	1:B:214:SER:HB2	2.56	0.41
1:B:27:GLN:O	1:B:30:LEU:CB	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:LEU:HD12	1:B:29:LEU:HA	1.80	0.41
1:A:59:ILE:HG13	1:A:60:HIS:N	2.25	0.41
1:A:125:TYR:HA	1:A:126:PRO:HA	1.79	0.40
1:B:103:HIS:HB2	3:B:1209:PLP:H2A3	2.03	0.40
1:B:123:THR:HB	1:B:129:ALA:H	1.86	0.40
1:A:208:HIS:CE1	3:A:1209:PLP:O1P	2.74	0.40
1:A:90:ALA:CB	1:A:115:LEU:HD21	2.51	0.40
1:A:97:VAL:O	1:A:97:VAL:HG23	2.21	0.40
1:B:63:LEU:HD21	1:B:219:VAL:HG22	2.03	0.40
1:A:190:ARG:NH1	4:A:1212:HOH:O	2.41	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%)	314 (92%)	19 (6%)	7 (2%)	5	7
1	B	340/371 (92%)	319 (94%)	17 (5%)	4 (1%)	10	16
All	All	680/742 (92%)	633 (93%)	36 (5%)	11 (2%)	7	11

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	349	TYR
1	A	215	GLY
1	A	249	ALA
1	B	215	GLY
1	A	224	GLU
1	B	250	ARG
1	A	31	GLU
1	A	176	TYR

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Mol	Chain	Res	Type
1	A	238	LYS
1	A	349	TYR
1	B	30	LEU

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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	701	-	4,4,4	0.36	0	6,6,6	0.17	0
2	SO4	A	601	-	4,4,4	0.42	0	6,6,6	0.11	0
2	SO4	A	602	-	4,4,4	0.36	0	6,6,6	0.18	0
2	SO4	B	702	-	4,4,4	0.39	0	6,6,6	0.31	0
3	PLP	B	1209	1	15,15,16	1.07	0	21,22,23	1.04	0
3	PLP	A	1209	1	15,15,16	1.08	0	21,22,23	1.00	1 (4%)

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	B	1209	1	-	5/6/6/8	0/1/1/1
3	PLP	A	1209	1	-	5/6/6/8	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	2.45	113.05	106.44

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1209	PLP	C5A-O4P-P-O1P
3	A	1209	PLP	C5A-O4P-P-O2P
3	A	1209	PLP	C5A-O4P-P-O3P
3	B	1209	PLP	C4-C5-C5A-O4P
3	B	1209	PLP	C6-C5-C5A-O4P
3	B	1209	PLP	C5A-O4P-P-O1P
3	B	1209	PLP	C5A-O4P-P-O2P
3	B	1209	PLP	C5A-O4P-P-O3P
3	A	1209	PLP	C6-C5-C5A-O4P
3	A	1209	PLP	C4-C5-C5A-O4P

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	602	SO4	1	0
2	B	702	SO4	1	0
3	B	1209	PLP	3	0
3	A	1209	PLP	2	0

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%)	0.33	15 (4%) 39 34	35, 57, 97, 123	0
1	B	344/371 (92%)	-0.06	14 (4%) 41 37	27, 44, 79, 104	0
All	All	688/742 (92%)	0.13	29 (4%) 40 36	27, 50, 93, 123	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	125	TYR	6.9
1	A	249	ALA	5.7
1	B	249	ALA	5.5
1	A	53	PRO	3.7
1	B	248	THR	3.6
1	B	53	PRO	3.5
1	A	32	TRP	3.4
1	A	228	GLU	3.1
1	A	9	PHE	2.8
1	B	9	PHE	2.8
1	B	250	ARG	2.7
1	B	332	LYS	2.7
1	B	215	GLY	2.6
1	A	30	LEU	2.5
1	A	250	ARG	2.4
1	B	317	ALA	2.4
1	A	103	HIS	2.4
1	A	248	THR	2.4
1	B	12	ILE	2.4
1	B	318	LYS	2.3
1	B	252	ALA	2.2
1	A	334	HIS	2.2
1	B	354	GLU	2.2
1	A	290	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	171	LYS	2.1
1	A	266	GLU	2.1
1	B	126	PRO	2.1
1	A	319	GLY	2.1
1	A	370	TYR	2.0

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	SO4	A	601	5/5	0.76	0.13	114,120,120,122	0
2	SO4	B	701	5/5	0.83	0.13	100,111,113,122	0
3	PLP	A	1209	15/16	0.92	0.15	46,92,107,110	0
3	PLP	B	1209	15/16	0.94	0.14	45,82,98,99	0
2	SO4	A	602	5/5	0.96	0.07	58,65,67,70	0
2	SO4	B	702	5/5	0.97	0.08	43,52,68,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.