



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:26 PM UTC

PDB ID : 2ZAU / pdb_00002zau
Title : Crystal structure of an N-terminally truncated selenophosphate synthetase from Aquifex aeolicus
Authors : Sekine, S.; Matsumoto, E.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-10-10
Resolution : 2.00 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 : 4-5-2 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

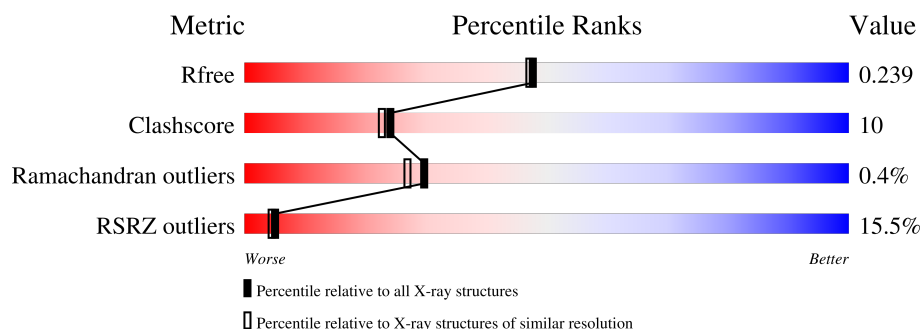
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	311	17% 78% 20% .
1	B	311	13% 77% 19% ..
1	C	311	16% 77% 20% .

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7533 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.

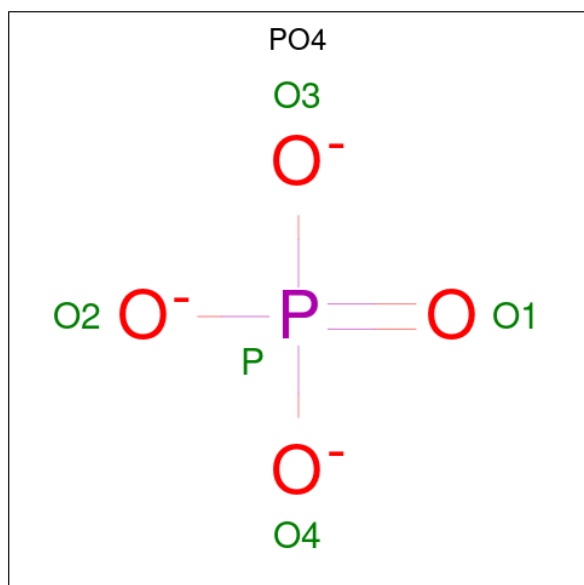
- Molecule 1 is a protein called Selenide, water dikinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2377	1530	383	456	8			
1	B	303	Total	C	N	O	S	0	0	0
			2348	1513	379	448	8			
1	C	304	Total	C	N	O	S	0	0	0
			2352	1517	380	447	8			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	MET	-	initiating methionine	UNP O67139
B	26	MET	-	initiating methionine	UNP O67139
C	26	MET	-	initiating methionine	UNP O67139

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

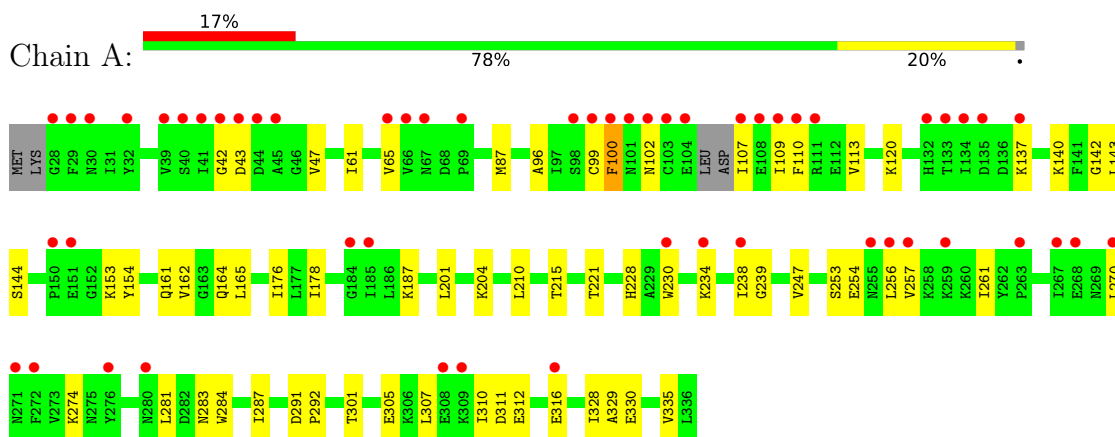
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	140	Total	O	0	0
			140	140		
3	B	162	Total	O	0	0
			162	162		
3	C	139	Total	O	0	0
			139	139		

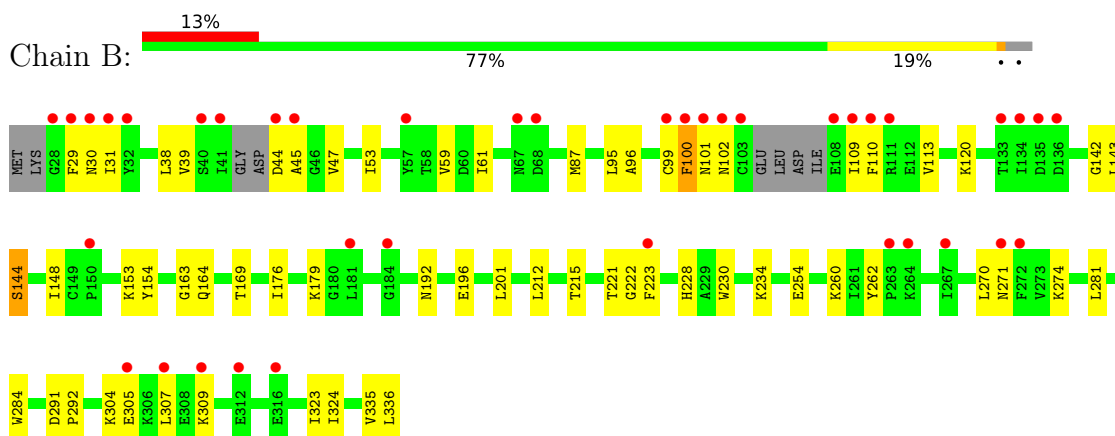
3 Residue-property plots [i](#)

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.

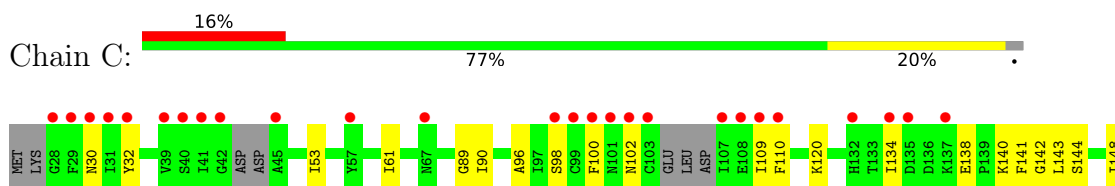
- Molecule 1: Selenide, water dikinase

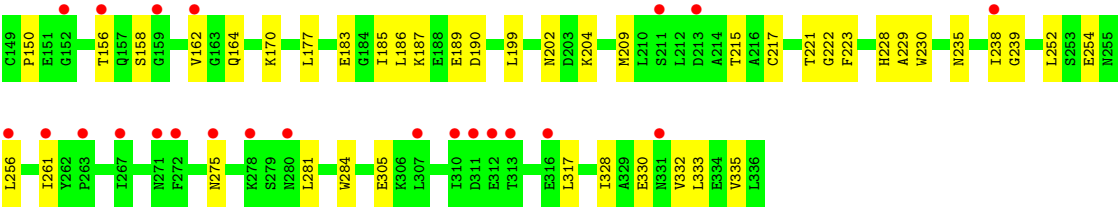


- Molecule 1: Selenide, water dikinase



- Molecule 1: Selenide, water dikinase





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	93.16Å 165.20Å 167.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.29 – 2.00 46.29 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.8 (46.29-2.00) 93.9 (46.29-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.42 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.212 , 0.239 0.212 , 0.239	Depositor DCC
R_{free} test set	4140 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.025 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.025 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7533	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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: PO4

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	0.38	0/2417	0.82	2/3277 (0.1%)
1	B	0.39	0/2387	0.84	4/3235 (0.1%)
1	C	0.36	0/2391	0.81	4/3240 (0.1%)
All	All	0.38	0/7195	0.82	10/9752 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	144	SER	N-CA-C	-8.84	94.84	109.07
1	A	144	SER	N-CA-C	-8.59	95.25	109.24
1	C	144	SER	N-CA-C	-7.08	97.66	109.07
1	B	324	ILE	N-CA-C	6.21	118.98	112.83
1	B	222	GLY	N-CA-C	5.87	121.51	113.99
1	C	222	GLY	N-CA-C	5.82	121.02	113.27
1	C	89	GLY	N-CA-C	5.74	120.75	112.82
1	A	247	VAL	N-CA-C	5.51	113.37	108.63
1	C	229	ALA	N-CA-C	-5.18	105.71	111.36
1	B	169	THR	N-CA-C	5.07	120.09	113.30

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	2377	0	2413	49	0
1	B	2348	0	2388	49	1
1	C	2352	0	2398	45	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
3	A	140	0	0	2	0
3	B	162	0	0	5	0
3	C	139	0	0	6	0
All	All	7533	0	7199	143	1

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 10.

All (143) 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:144:SER:HB3	3:B:1412:HOH:O	1.56	1.05
1:B:95:LEU:HB2	3:B:1412:HOH:O	1.68	0.92
1:A:162:VAL:HG22	1:A:238:ILE:HD11	1.53	0.89
1:C:134:ILE:H	1:C:134:ILE:HD13	1.49	0.78
1:B:44:ASP:HB3	1:B:59:VAL:HG21	1.67	0.75
1:A:254:GLU:HG3	1:A:284:TRP:CZ3	2.27	0.70
1:B:87:MET:HE2	1:B:154:TYR:HE1	1.59	0.68
1:A:238:ILE:HG13	1:A:328:ILE:O	1.93	0.68
1:A:253:SER:O	1:A:257:VAL:HG22	1.95	0.66
1:A:312:GLU:O	1:A:316:GLU:HG3	1.96	0.66
1:A:176:ILE:HD13	1:A:257:VAL:HG13	1.76	0.66
1:A:87:MET:HE2	1:A:154:TYR:HE1	1.60	0.65
1:C:330:GLU:HG2	1:C:332:VAL:HG22	1.80	0.63
1:C:221:THR:HG21	3:C:1027:HOH:O	1.99	0.61
1:C:223:PHE:HB2	1:C:228:HIS:CE1	2.36	0.61
1:A:120:LYS:HD2	1:A:201:LEU:HD13	1.83	0.61
1:B:221:THR:H	1:B:228:HIS:CE1	2.19	0.60
1:B:87:MET:HE2	1:B:154:TYR:CE1	2.35	0.60
1:B:254:GLU:HG3	1:B:284:TRP:CZ3	2.38	0.59
1:C:183:GLU:HB3	1:C:185:ILE:HG13	1.84	0.59
1:A:47:VAL:HG21	1:A:154:TYR:OH	2.03	0.59
1:A:162:VAL:CG2	1:A:238:ILE:HD11	2.30	0.58
1:A:270:LEU:O	1:A:274:LYS:HG2	2.04	0.58
1:C:162:VAL:HB	1:C:238:ILE:HD11	1.86	0.58
1:B:120:LYS:HD2	1:B:201:LEU:HD13	1.86	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ILE:CD1	1:A:257:VAL:HG13	2.34	0.57
1:A:161:GLN:HG2	3:A:1274:HOH:O	2.04	0.57
1:C:156:THR:HG23	1:C:158:SER:H	1.69	0.57
1:C:185:ILE:HD13	1:C:261:ILE:HD13	1.87	0.57
1:A:238:ILE:HG13	1:A:239:GLY:H	1.68	0.56
1:B:305:GLU:H	1:B:305:GLU:CD	2.13	0.56
1:A:153:LYS:HD3	1:A:210:LEU:HD22	1.87	0.56
1:B:53:ILE:HD11	1:B:148:ILE:HG23	1.87	0.55
1:C:177:LEU:HD21	1:C:252:LEU:HB3	1.87	0.55
1:B:153:LYS:HB3	3:B:1387:HOH:O	2.06	0.55
1:B:53:ILE:HD11	1:B:148:ILE:CG2	2.36	0.55
1:C:238:ILE:HG13	1:C:239:GLY:H	1.70	0.55
1:A:87:MET:HE2	1:A:154:TYR:CE1	2.42	0.54
1:A:238:ILE:HD11	1:A:329:ALA:HA	1.89	0.54
1:B:192:ASN:O	1:B:196:GLU:HG3	2.07	0.54
1:C:223:PHE:H	1:C:228:HIS:HE1	1.55	0.54
1:A:61:ILE:HA	1:A:142:GLY:HA3	1.89	0.54
1:C:238:ILE:HG13	1:C:328:ILE:O	2.08	0.53
1:B:260:LYS:H	1:B:260:LYS:HE2	1.74	0.53
1:A:65:VAL:HG13	1:A:178:ILE:CD1	2.40	0.52
1:A:65:VAL:HG13	1:A:178:ILE:HD13	1.92	0.52
1:B:179:LYS:HD2	1:B:262:TYR:C	2.35	0.52
1:B:45:ALA:HB2	3:B:1383:HOH:O	2.09	0.51
1:A:270:LEU:HD22	1:A:274:LYS:HD3	1.92	0.51
1:A:239:GLY:HA3	1:A:330:GLU:O	2.10	0.51
1:B:192:ASN:H	1:B:192:ASN:HD22	1.58	0.51
1:B:29:PHE:CZ	1:B:31:ILE:HD11	2.46	0.51
1:C:158:SER:HB3	1:C:235:ASN:HB3	1.91	0.51
1:C:134:ILE:H	1:C:134:ILE:CD1	2.23	0.50
1:A:305:GLU:CD	1:A:305:GLU:H	2.19	0.50
1:C:53:ILE:HD12	3:C:1399:HOH:O	2.12	0.50
1:C:90:ILE:HD12	1:C:150:PRO:HG2	1.94	0.50
1:A:102:ASN:O	1:A:137:LYS:HA	2.12	0.49
1:B:44:ASP:HB3	1:B:59:VAL:CG2	2.38	0.49
1:C:138:GLU:HB3	1:C:140:LYS:NZ	2.28	0.49
1:B:29:PHE:O	1:B:30:ASN:ND2	2.43	0.49
1:B:223:PHE:H	1:B:228:HIS:CE1	2.31	0.49
1:A:140:LYS:NZ	1:A:140:LYS:HB2	2.28	0.48
1:B:260:LYS:HE2	1:B:260:LYS:N	2.27	0.48
1:C:156:THR:HG22	3:C:1340:HOH:O	2.12	0.48
1:C:185:ILE:CD1	1:C:261:ILE:HD13	2.43	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:TRP:HE1	1:A:234:LYS:HD3	1.77	0.48
1:A:164:GLN:HE22	1:A:215:THR:CG2	2.26	0.48
1:C:100:PHE:C	1:C:102:ASN:H	2.20	0.48
1:C:61:ILE:HA	1:C:142:GLY:HA3	1.96	0.48
1:B:61:ILE:HA	1:B:142:GLY:HA3	1.96	0.47
1:B:230:TRP:NE1	1:B:234:LYS:HD3	2.29	0.47
1:C:187:LYS:HG2	1:C:190:ASP:OD2	2.15	0.47
1:C:228:HIS:HD2	3:C:1343:HOH:O	1.97	0.46
1:C:96:ALA:HA	1:C:143:LEU:HD23	1.98	0.46
1:A:100:PHE:C	1:A:102:ASN:H	2.23	0.46
1:A:109:ILE:HG13	1:A:110:PHE:N	2.31	0.46
1:B:95:LEU:HD12	3:B:1412:HOH:O	2.16	0.46
1:C:170:LYS:HE3	1:C:202:ASN:OD1	2.16	0.46
1:B:212:LEU:HD23	1:B:309:LYS:HE3	1.97	0.46
1:C:186:LEU:HD11	1:C:256:LEU:HD11	1.96	0.46
1:A:99:CYS:O	1:A:100:PHE:O	2.34	0.46
1:A:281:LEU:HD12	1:A:335:VAL:HG12	1.98	0.46
1:B:307:LEU:HD11	1:B:323:ILE:CD1	2.46	0.46
1:A:283:ASN:O	1:A:287:ILE:HG12	2.16	0.45
1:B:223:PHE:H	1:B:228:HIS:HE1	1.62	0.45
1:C:199:LEU:HG	3:C:1128:HOH:O	2.16	0.45
1:A:204:LYS:HE2	3:A:1192:HOH:O	2.17	0.45
1:A:270:LEU:HD12	1:A:287:ILE:HD13	1.99	0.45
1:C:238:ILE:HG23	1:C:239:GLY:N	2.30	0.45
1:C:254:GLU:HG3	1:C:284:TRP:CZ3	2.51	0.45
1:B:270:LEU:O	1:B:274:LYS:HG3	2.17	0.44
1:C:98:SER:HB3	1:C:141:PHE:HD1	1.81	0.44
1:A:310:ILE:HG23	1:A:311:ASP:N	2.32	0.44
1:B:100:PHE:O	1:B:101:ASN:HB3	2.17	0.44
1:B:100:PHE:C	1:B:102:ASN:H	2.26	0.44
1:C:230:TRP:HB2	1:C:333:LEU:HD11	1.99	0.44
1:B:164:GLN:HE22	1:B:215:THR:CG2	2.29	0.44
1:C:30:ASN:HB3	1:C:32:TYR:CE1	2.53	0.44
1:B:47:VAL:HG21	1:B:154:TYR:OH	2.18	0.44
1:C:221:THR:H	1:C:228:HIS:CE1	2.34	0.44
1:C:164:GLN:HE22	1:C:215:THR:CG2	2.30	0.44
1:B:230:TRP:HE1	1:B:234:LYS:HD3	1.83	0.44
1:A:256:LEU:HB3	1:A:261:ILE:HB	1.99	0.43
1:A:281:LEU:CD1	1:A:335:VAL:HG12	2.48	0.43
1:A:230:TRP:NE1	1:A:234:LYS:HD3	2.34	0.43
1:C:164:GLN:HE22	1:C:215:THR:HG21	1.84	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:LYS:HD2	1:C:189:GLU:HB2	2.00	0.43
1:C:209:MET:HE2	1:C:217:CYS:SG	2.59	0.43
1:B:99:CYS:O	1:B:100:PHE:C	2.62	0.43
1:A:96:ALA:HA	1:A:143:LEU:HD23	2.01	0.43
1:A:162:VAL:HG22	1:A:238:ILE:CD1	2.36	0.43
1:B:109:ILE:O	1:B:113:VAL:HG23	2.18	0.42
1:B:109:ILE:HG13	1:B:110:PHE:N	2.34	0.42
1:B:38:LEU:HD23	1:B:39:VAL:N	2.34	0.42
1:B:291:ASP:HA	1:B:292:PRO:HD3	1.90	0.42
1:A:187:LYS:HB2	1:A:187:LYS:NZ	2.34	0.42
1:B:223:PHE:HB2	1:B:228:HIS:CE1	2.54	0.42
1:C:281:LEU:HD12	1:C:335:VAL:HG12	2.02	0.42
1:B:291:ASP:OD2	1:B:292:PRO:HD2	2.20	0.42
1:B:38:LEU:HD23	1:B:38:LEU:C	2.45	0.42
1:B:29:PHE:C	1:B:30:ASN:HD22	2.28	0.41
1:C:177:LEU:CD2	1:C:252:LEU:HB3	2.50	0.41
1:B:176:ILE:HG13	1:B:291:ASP:CG	2.45	0.41
1:A:221:THR:H	1:A:228:HIS:CD2	2.38	0.41
1:B:163:GLY:O	1:B:304:LYS:HE2	2.20	0.41
1:C:120:LYS:HA	1:C:120:LYS:HE2	2.01	0.41
1:C:109:ILE:HG13	1:C:110:PHE:N	2.36	0.41
1:B:336:LEU:HD23	1:B:336:LEU:N	2.35	0.41
1:A:291:ASP:HA	1:A:292:PRO:HD3	1.96	0.41
1:B:96:ALA:HA	1:B:143:LEU:HD23	2.03	0.41
1:C:275:ASN:ND2	3:C:1359:HOH:O	2.53	0.41
1:A:165:LEU:O	1:A:301:THR:HA	2.21	0.41
1:A:165:LEU:HD11	1:A:307:LEU:HD11	2.03	0.41
1:C:204:LYS:HD3	1:C:317:LEU:O	2.21	0.41
1:C:305:GLU:OE2	1:C:305:GLU:N	2.52	0.41
1:A:61:ILE:HG22	1:A:142:GLY:HA3	2.03	0.40
1:A:238:ILE:HD12	1:A:238:ILE:HA	1.89	0.40
1:C:148:ILE:O	1:C:150:PRO:HD3	2.20	0.40
1:A:109:ILE:O	1:A:113:VAL:HG23	2.21	0.40
1:A:107:ILE:O	1:A:107:ILE:HG22	2.21	0.40
1:B:234:LYS:HB2	1:B:234:LYS:NZ	2.37	0.40
1:B:281:LEU:HD12	1:B:335:VAL:HG12	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:ASN:OD1	1:B:271:ASN:OD1[4_576]	1.75	0.45

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	303/311 (97%)	292 (96%)	8 (3%)	3 (1%)	12	8
1	B	297/311 (96%)	286 (96%)	10 (3%)	1 (0%)	36	35
1	C	298/311 (96%)	287 (96%)	11 (4%)	0	100	100
All	All	898/933 (96%)	865 (96%)	29 (3%)	4 (0%)	30	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	ASP
1	A	100	PHE
1	B	100	PHE
1	A	42	GLY

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](#)

3 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	PO4	B	903	-	4,4,4	2.00	3 (75%)	6,6,6	0.49	0
2	PO4	C	901	-	4,4,4	1.75	1 (25%)	6,6,6	0.48	0
2	PO4	A	902	-	4,4,4	1.80	3 (75%)	6,6,6	0.46	0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	903	PO4	P-O2	-2.29	1.48	1.54
2	B	903	PO4	P-O4	-2.15	1.48	1.54
2	B	903	PO4	P-O3	-2.07	1.48	1.54
2	A	902	PO4	P-O2	-2.03	1.48	1.54
2	A	902	PO4	P-O3	-2.03	1.48	1.54
2	A	902	PO4	P-O4	-2.02	1.48	1.54
2	C	901	PO4	P-O4	-2.01	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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	307/311 (98%)	0.84	54 (17%) 4 3	21, 35, 67, 95	0
1	B	303/311 (97%)	0.56	39 (12%) 7 6	18, 33, 56, 78	0
1	C	304/311 (97%)	0.91	49 (16%) 4 4	21, 37, 65, 91	0
All	All	914/933 (97%)	0.77	142 (15%) 5 4	18, 35, 65, 95	0

All (142) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	107	ILE	8.6
1	C	100	PHE	7.3
1	C	107	ILE	7.0
1	B	100	PHE	6.7
1	C	99	CYS	6.5
1	B	109	ILE	6.2
1	C	103	CYS	6.2
1	A	99	CYS	5.6
1	C	28	GLY	5.6
1	A	100	PHE	5.4
1	A	272	PHE	5.3
1	A	28	GLY	5.3
1	B	45	ALA	5.1
1	A	134	ILE	4.7
1	A	41	ILE	4.4
1	C	41	ILE	4.4
1	A	42	GLY	4.2
1	C	267	ILE	4.1
1	B	135	ASP	4.1
1	C	271	ASN	4.1
1	A	267	ILE	4.1
1	B	99	CYS	4.1
1	C	272	PHE	4.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	109	ILE	3.9
1	A	30	ASN	3.9
1	A	102	ASN	3.9
1	B	101	ASN	3.8
1	C	134	ILE	3.7
1	B	271	ASN	3.7
1	B	41	ILE	3.7
1	A	110	PHE	3.7
1	A	109	ILE	3.6
1	C	110	PHE	3.6
1	A	29	PHE	3.5
1	A	255	ASN	3.5
1	C	45	ALA	3.5
1	C	67	ASN	3.4
1	C	40	SER	3.4
1	C	316	GLU	3.4
1	A	66	VAL	3.4
1	A	238	ILE	3.4
1	B	103	CYS	3.4
1	B	110	PHE	3.4
1	B	28	GLY	3.3
1	C	238	ILE	3.3
1	C	312	GLU	3.3
1	B	184	GLY	3.2
1	A	45	ALA	3.2
1	B	31	ILE	3.2
1	A	104	GLU	3.2
1	C	135	ASP	3.1
1	A	263	PRO	3.1
1	B	67	ASN	3.1
1	C	101	ASN	3.1
1	A	270	LEU	3.1
1	B	272	PHE	3.1
1	A	101	ASN	3.0
1	B	44	ASP	3.0
1	A	67	ASN	3.0
1	C	98	SER	2.9
1	C	42	GLY	2.9
1	A	108	GLU	2.9
1	B	316	GLU	2.9
1	C	102	ASN	2.9
1	A	44	ASP	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	263	PRO	2.9
1	A	40	SER	2.8
1	C	30	ASN	2.8
1	B	32	TYR	2.8
1	A	111	ARG	2.8
1	C	137	LYS	2.8
1	B	108	GLU	2.8
1	B	307	LEU	2.8
1	A	185	ILE	2.8
1	C	310	ILE	2.8
1	A	43	ASP	2.8
1	A	133	THR	2.8
1	C	31	ILE	2.8
1	A	271	ASN	2.7
1	C	263	PRO	2.7
1	C	162	VAL	2.7
1	B	181	LEU	2.7
1	A	98	SER	2.7
1	B	102	ASN	2.6
1	C	311	ASP	2.6
1	A	103	CYS	2.6
1	A	184	GLY	2.6
1	A	257	VAL	2.6
1	C	156	THR	2.6
1	B	111	ARG	2.6
1	C	331	ASN	2.6
1	C	108	GLU	2.5
1	B	264	LYS	2.5
1	A	135	ASP	2.5
1	C	132	HIS	2.5
1	A	65	VAL	2.5
1	B	305	GLU	2.5
1	A	132	HIS	2.5
1	C	159	GLY	2.5
1	B	134	ILE	2.5
1	A	308	GLU	2.4
1	C	152	GLY	2.4
1	C	280	ASN	2.4
1	B	133	THR	2.4
1	B	312	GLU	2.4
1	A	234	LYS	2.4
1	B	309	LYS	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	32	TYR	2.4
1	A	280	ASN	2.4
1	C	275	ASN	2.4
1	C	32	TYR	2.3
1	C	29	PHE	2.3
1	A	309	LYS	2.3
1	A	151	GLU	2.3
1	B	30	ASN	2.3
1	A	256	LEU	2.3
1	B	136	ASP	2.3
1	C	307	LEU	2.3
1	A	39	VAL	2.2
1	A	150	PRO	2.2
1	C	213	ASP	2.2
1	A	276	TYR	2.2
1	B	57	TYR	2.2
1	A	137	LYS	2.2
1	B	150	PRO	2.2
1	B	68	ASP	2.2
1	B	29	PHE	2.2
1	C	313	THR	2.2
1	B	40	SER	2.2
1	A	230	TRP	2.2
1	B	223	PHE	2.1
1	B	267	ILE	2.1
1	A	268	GLU	2.1
1	C	39	VAL	2.1
1	C	278	LYS	2.1
1	A	316	GLU	2.0
1	C	211	SER	2.1
1	A	69	PRO	2.0
1	C	261	ILE	2.0
1	C	256	LEU	2.0
1	A	259	LYS	2.0
1	C	57	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [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	PO4	B	903	5/5	0.83	0.17	60,62,64,64	0
2	PO4	C	901	5/5	0.93	0.11	50,51,54,54	0
2	PO4	A	902	5/5	0.95	0.10	54,55,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.