



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:14 PM UTC

PDB ID : 3U0O / pdb_00003u0o
Title : The crystal structure of selenophosphate synthetase from E. coli
Authors : Noinaj, N.; Wattanasak, R.; Wally, J.; Stadtman, T.; Buchanan, S.K.
Deposited on : 2011-09-28
Resolution : 2.25 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
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

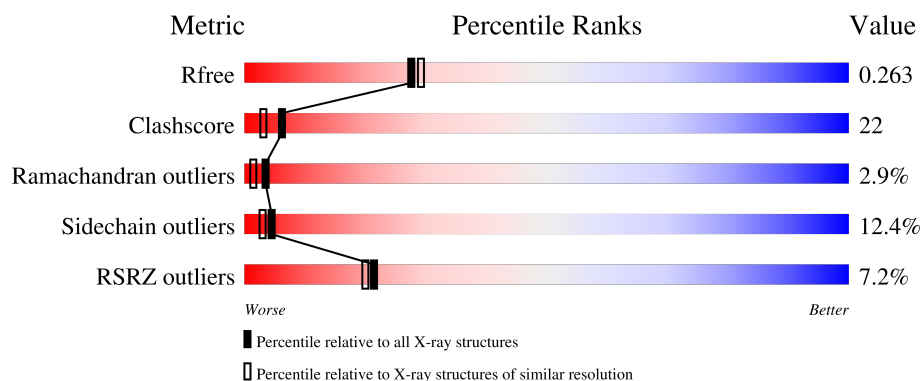
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	347	5% 68% 19% 5% 8%
1	B	347	8% 61% 26% 8% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4621 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	319	Total	C	N	O	S	0	0	0
			2209	1391	374	428	16			
1	B	335	Total	C	N	O	S	0	0	0
			2358	1493	396	451	18			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

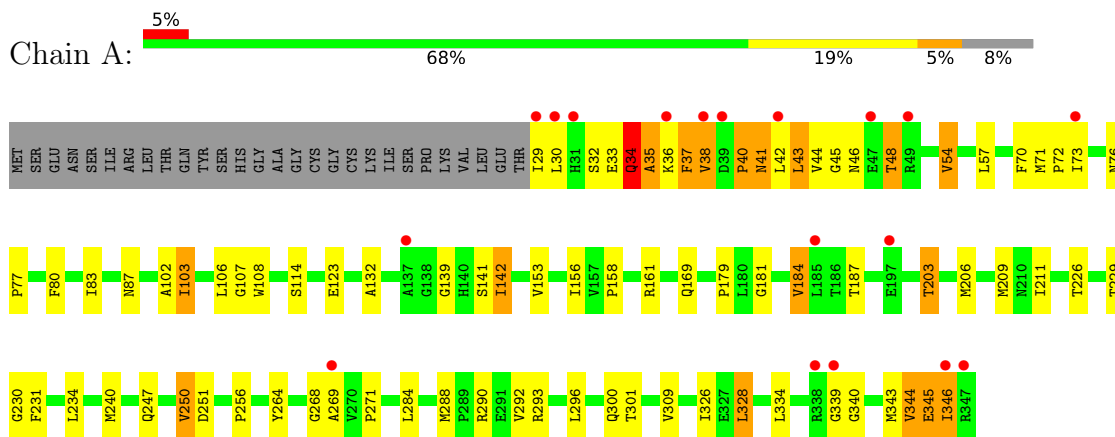
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	31	Total	O	0	0
			31	31		
3	B	21	Total	O	0	0
			21	21		

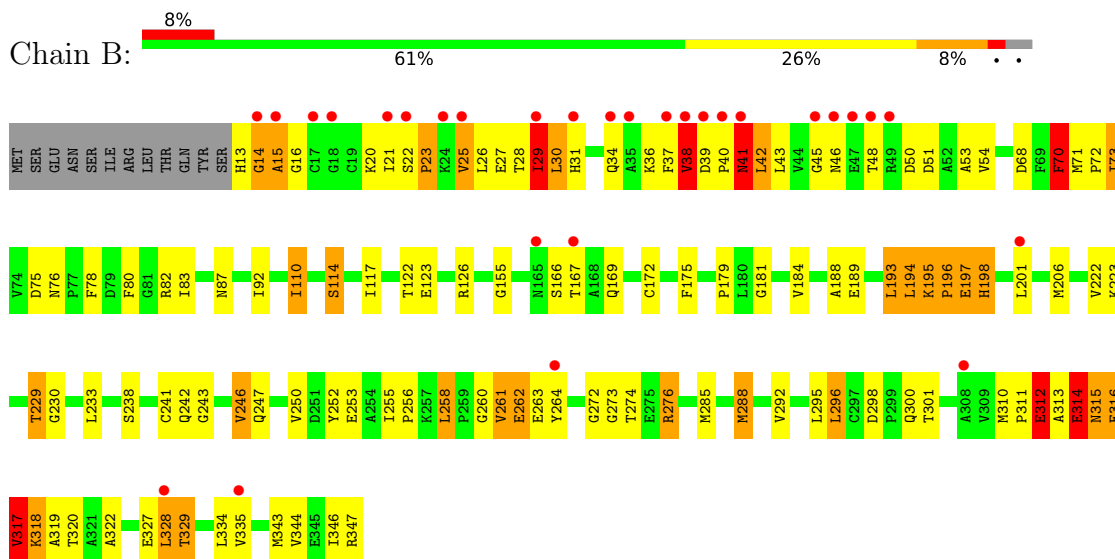
3 Residue-property plots [i](#)

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.35Å 102.35Å 140.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.98 – 2.25 28.98 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.8 (28.98-2.25) 99.8 (28.98-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 2.24Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_865)	Depositor
R, R_{free}	0.222 , 0.264 0.221 , 0.263	Depositor DCC
R_{free} test set	2059 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	57.3	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 79.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4621	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: MG

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	0.65	1/2247 (0.0%)	1.03	9/3064 (0.3%)
1	B	0.73	1/2401 (0.0%)	1.16	21/3273 (0.6%)
All	All	0.69	2/4648 (0.0%)	1.10	30/6337 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	344	VAL	CA-CB	-6.82	1.46	1.54
1	B	195	LYS	C-O	-5.51	1.19	1.24

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	15	ALA	N-CA-C	-15.48	95.08	112.57
1	B	29	ILE	N-CA-C	-8.02	98.06	110.09
1	A	231	PHE	N-CA-C	-7.58	104.03	113.28
1	B	243	GLY	N-CA-C	-6.64	104.04	112.54
1	B	41	ASN	N-CA-C	6.59	120.09	109.02
1	B	39	ASP	N-CA-C	6.36	117.92	108.76
1	B	114	SER	CA-C-N	6.16	125.84	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	114	SER	C-N-CA	6.16	125.84	119.56
1	A	345	GLU	N-CA-C	-6.13	97.75	110.80
1	A	114	SER	CA-C-N	5.73	126.12	119.47
1	A	114	SER	C-N-CA	5.73	126.12	119.47
1	A	57	LEU	N-CA-C	-5.72	101.17	110.20
1	B	250	VAL	N-CA-C	5.61	115.96	108.11
1	B	315	ASN	CA-C-N	-5.60	112.28	121.19
1	B	315	ASN	C-N-CA	-5.60	112.28	121.19
1	A	71	MET	N-CA-C	-5.57	102.89	110.36
1	B	70	PHE	N-CA-C	5.53	116.37	108.74
1	B	198	HIS	N-CA-C	-5.36	106.61	112.72
1	A	340	GLY	N-CA-C	5.32	118.05	111.93
1	B	329	THR	N-CA-C	5.31	117.06	108.41
1	A	34	GLN	N-CA-C	-5.25	99.61	110.80
1	A	309	VAL	N-CA-C	5.23	115.51	107.77
1	B	76	ASN	CA-C-N	5.11	124.73	119.05
1	B	76	ASN	C-N-CA	5.11	124.73	119.05
1	B	36	LYS	CB-CA-C	-5.10	110.71	116.63
1	B	316	GLU	CA-C-N	-5.10	112.78	121.97
1	B	316	GLU	C-N-CA	-5.10	112.78	121.97
1	B	229	THR	CB-CA-C	-5.07	103.22	110.62
1	B	328	LEU	CA-C-N	-5.01	115.94	123.00
1	B	328	LEU	C-N-CA	-5.01	115.94	123.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	45	GLY	Peptide
1	A	48	THR	Peptide
1	B	22	SER	Peptide
1	B	29	ILE	Peptide
1	B	41	ASN	Peptide

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	2209	0	2091	69	0
1	B	2358	0	2270	137	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	31	0	0	0	0
3	B	21	0	0	1	0
All	All	4621	0	4361	200	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 22.

All (200) 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:34:GLN:HA	1:A:35:ALA:CB	1.51	1.34
1:B:194:LEU:HD13	1:B:196:PRO:CD	1.60	1.32
1:A:37:PHE:O	1:A:38:VAL:HG22	1.28	1.28
1:B:194:LEU:C	1:B:194:LEU:HD12	1.56	1.26
1:B:194:LEU:HD12	1:B:194:LEU:O	1.35	1.26
1:B:314:GLU:O	1:B:318:LYS:HG3	1.28	1.25
1:B:311:PRO:O	1:B:313:ALA:N	1.82	1.12
1:A:37:PHE:O	1:A:38:VAL:CG2	1.97	1.11
1:A:34:GLN:HA	1:A:35:ALA:HB2	1.18	1.10
1:A:34:GLN:HA	1:A:35:ALA:HB3	1.33	1.10
1:A:41:ASN:HB2	1:A:42:LEU:HA	1.10	1.10
1:B:317:VAL:HG12	1:B:318:LYS:N	1.53	1.09
1:B:194:LEU:HD13	1:B:196:PRO:HD2	1.18	1.08
1:A:34:GLN:CA	1:A:35:ALA:CB	2.30	1.07
1:A:37:PHE:C	1:A:38:VAL:HG22	1.75	1.06
1:A:37:PHE:O	1:A:38:VAL:HG13	1.56	1.05
1:B:194:LEU:C	1:B:194:LEU:CD1	2.30	1.05
1:B:260:GLY:O	1:B:264:TYR:HD1	1.37	1.04
1:A:34:GLN:CA	1:A:35:ALA:HB2	1.85	1.03
1:B:193:LEU:HD22	1:B:194:LEU:N	1.74	1.03
1:B:314:GLU:O	1:B:318:LYS:CG	2.13	0.97
1:A:33:GLU:O	1:A:34:GLN:O	1.83	0.97
1:B:315:ASN:O	1:B:316:GLU:C	2.04	0.94
1:B:193:LEU:CD2	1:B:194:LEU:N	2.30	0.94
1:B:313:ALA:O	1:B:315:ASN:N	2.01	0.94
1:A:41:ASN:HB2	1:A:42:LEU:CA	1.99	0.92
1:B:311:PRO:C	1:B:313:ALA:H	1.75	0.92
1:B:194:LEU:CD1	1:B:196:PRO:CD	2.48	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:PHE:O	1:A:38:VAL:CG1	2.16	0.91
1:A:41:ASN:CB	1:A:42:LEU:HA	1.99	0.91
1:B:260:GLY:O	1:B:264:TYR:CD1	2.25	0.90
1:A:36:LYS:O	1:A:37:PHE:CB	2.21	0.89
1:B:252:TYR:O	1:B:252:TYR:CD2	2.30	0.85
1:B:194:LEU:CD1	1:B:196:PRO:HD2	2.06	0.85
1:A:103:ILE:HG23	1:A:139:GLY:HA2	1.57	0.85
1:B:194:LEU:HD13	1:B:196:PRO:CG	2.06	0.84
1:B:252:TYR:CE1	1:B:292:VAL:HG13	2.13	0.84
1:B:316:GLU:O	1:B:317:VAL:O	1.97	0.83
1:B:193:LEU:CD2	1:B:194:LEU:H	1.92	0.81
1:B:317:VAL:HG12	1:B:318:LYS:H	1.40	0.81
1:A:37:PHE:O	1:A:38:VAL:CB	2.29	0.81
1:B:317:VAL:CG1	1:B:318:LYS:N	2.30	0.81
1:B:41:ASN:CG	1:B:42:LEU:H	1.89	0.80
1:B:193:LEU:HD22	1:B:193:LEU:C	2.04	0.80
1:A:37:PHE:C	1:A:38:VAL:CG2	2.49	0.80
1:B:29:ILE:HG13	1:B:276:ARG:HH21	1.46	0.80
1:B:194:LEU:CD1	1:B:196:PRO:N	2.46	0.79
1:A:251:ASP:HB2	1:A:345:GLU:OE2	1.82	0.79
1:B:193:LEU:CD2	1:B:193:LEU:C	2.55	0.77
1:B:316:GLU:O	1:B:317:VAL:C	2.20	0.77
1:B:313:ALA:O	1:B:314:GLU:C	2.30	0.73
1:A:251:ASP:CB	1:A:345:GLU:OE2	2.36	0.73
1:B:194:LEU:HD13	1:B:196:PRO:N	2.05	0.72
1:B:327:GLU:HA	1:B:328:LEU:HB2	1.71	0.72
1:B:193:LEU:HD23	1:B:194:LEU:H	1.54	0.71
1:B:14:GLY:C	1:B:16:GLY:H	1.96	0.71
1:B:196:PRO:C	1:B:198:HIS:H	1.99	0.71
1:B:252:TYR:O	1:B:252:TYR:CG	2.44	0.70
1:B:41:ASN:CG	1:B:42:LEU:N	2.50	0.69
1:B:253:GLU:C	1:B:255:ILE:H	2.01	0.69
1:A:344:VAL:C	1:A:346:ILE:N	2.49	0.68
1:B:15:ALA:CB	1:B:75:ASP:OD1	2.41	0.68
1:B:193:LEU:HD23	1:B:194:LEU:N	2.09	0.67
1:A:142:ILE:HD13	1:B:71:MET:HE1	1.75	0.67
1:B:188:ALA:HB1	1:B:264:TYR:CD2	2.29	0.67
1:B:311:PRO:C	1:B:313:ALA:N	2.38	0.67
1:A:73:ILE:HD11	1:A:83:ILE:HG21	1.76	0.66
1:B:313:ALA:C	1:B:315:ASN:N	2.53	0.66
1:A:344:VAL:O	1:A:346:ILE:N	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:VAL:O	1:B:319:ALA:N	2.29	0.66
1:B:83:ILE:HG21	1:B:206:MET:HE2	1.77	0.65
1:B:196:PRO:O	1:B:198:HIS:N	2.29	0.65
1:B:318:LYS:O	1:B:319:ALA:C	2.38	0.65
1:B:315:ASN:O	1:B:316:GLU:O	2.15	0.65
1:B:317:VAL:O	1:B:320:THR:N	2.30	0.64
1:A:41:ASN:HA	1:A:43:LEU:H	1.64	0.63
1:A:44:VAL:HB	1:A:54:VAL:HG22	1.81	0.62
1:B:41:ASN:OD1	1:B:42:LEU:N	2.32	0.62
1:A:33:GLU:C	1:A:34:GLN:O	2.42	0.62
1:B:38:VAL:HG22	1:B:40:PRO:HD3	1.81	0.62
1:B:188:ALA:CB	1:B:264:TYR:CD2	2.83	0.62
1:B:317:VAL:O	1:B:318:LYS:C	2.42	0.62
1:B:194:LEU:HD11	1:B:196:PRO:HB2	1.82	0.62
1:B:194:LEU:HG	1:B:264:TYR:OH	2.00	0.61
1:A:108:TRP:HB3	1:B:31:HIS:NE2	2.15	0.61
1:B:123:GLU:OE2	1:B:126:ARG:NH1	2.34	0.60
1:A:158:PRO:HG2	1:A:161:ARG:HD3	1.83	0.60
1:B:68:ASP:OD2	1:B:87:ASN:ND2	2.34	0.60
1:A:203:THR:HA	1:A:206:MET:HE3	1.84	0.60
1:B:230:GLY:H	1:B:300:GLN:H	1.50	0.60
1:B:41:ASN:C	1:B:43:LEU:H	2.09	0.59
1:B:311:PRO:O	1:B:312:GLU:C	2.42	0.59
1:A:37:PHE:C	1:A:38:VAL:HG13	2.25	0.59
1:B:23:PRO:HB3	1:B:273:GLY:HA2	1.85	0.59
1:B:70:PHE:HE1	1:B:73:ILE:HD11	1.66	0.59
1:B:78:PHE:CE2	1:B:82:ARG:HD2	2.37	0.59
1:B:193:LEU:HD11	1:B:198:HIS:O	2.03	0.58
1:A:103:ILE:CG2	1:A:139:GLY:HA2	2.33	0.58
1:A:247:GLN:NE2	1:A:339:GLY:O	2.36	0.58
1:B:195:LYS:N	1:B:196:PRO:HD2	2.19	0.57
1:A:344:VAL:C	1:A:346:ILE:H	2.09	0.57
1:B:194:LEU:O	1:B:194:LEU:CD1	2.30	0.57
1:A:268:GLY:HA2	1:A:269:ALA:HB3	1.86	0.56
1:B:181:GLY:HA3	1:B:261:VAL:HG11	1.87	0.56
1:A:229:THR:HA	1:A:300:GLN:O	2.06	0.56
1:B:194:LEU:CD1	1:B:196:PRO:CG	2.83	0.56
1:B:189:GLU:HA	1:B:193:LEU:HB2	1.87	0.55
1:A:34:GLN:CA	1:A:35:ALA:HB3	2.17	0.54
1:A:234:LEU:HD21	1:A:296:LEU:HD23	1.89	0.54
1:A:33:GLU:HG3	1:B:122:THR:HG21	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:LYS:O	1:B:25:VAL:HG21	2.08	0.54
1:B:194:LEU:HG	1:B:264:TYR:HH	1.72	0.54
1:B:256:PRO:HD3	1:B:329:THR:HG21	1.90	0.53
1:B:92:ILE:HD13	1:B:155:GLY:HA3	1.89	0.53
1:B:15:ALA:HB2	1:B:75:ASP:OD1	2.08	0.53
1:A:132:ALA:HB2	1:A:209:MET:HE3	1.91	0.53
1:A:103:ILE:HD11	1:B:53:ALA:HB3	1.91	0.53
1:B:252:TYR:CD2	1:B:252:TYR:C	2.87	0.52
1:B:258:LEU:O	1:B:261:VAL:HG22	2.10	0.52
1:A:33:GLU:N	1:A:33:GLU:OE1	2.43	0.51
1:B:38:VAL:HG13	1:B:40:PRO:HD2	1.91	0.51
1:B:37:PHE:O	1:B:38:VAL:HB	2.10	0.51
1:B:117:ILE:HD12	1:B:117:ILE:H	1.73	0.51
1:A:187:THR:HG21	1:A:271:PRO:HB3	1.92	0.51
1:B:45:GLY:HA3	1:B:54:VAL:HB	1.92	0.51
1:A:73:ILE:HD13	1:A:301:THR:HG21	1.93	0.50
1:A:250:VAL:CG2	1:A:344:VAL:HG13	2.41	0.50
1:B:195:LYS:O	1:B:196:PRO:O	2.29	0.50
1:A:40:PRO:O	1:A:41:ASN:O	2.29	0.50
1:B:288:MET:HE3	1:B:346:ILE:HD12	1.93	0.50
1:B:233:LEU:HD23	1:B:296:LEU:HD12	1.94	0.50
1:A:345:GLU:O	1:A:346:ILE:O	2.30	0.49
1:B:263:GLU:OE2	1:B:263:GLU:N	2.37	0.49
1:A:70:PHE:HZ	1:A:87:ASN:ND2	2.10	0.49
1:B:15:ALA:HB3	1:B:75:ASP:OD1	2.10	0.49
1:A:234:LEU:HD11	1:A:296:LEU:HB3	1.94	0.49
1:B:188:ALA:O	1:B:193:LEU:HB2	2.12	0.49
1:A:264:TYR:O	1:A:268:GLY:HA2	2.13	0.49
1:B:195:LYS:N	1:B:196:PRO:CD	2.76	0.48
1:B:313:ALA:C	1:B:315:ASN:H	2.20	0.48
1:B:29:ILE:H	1:B:276:ARG:NH2	2.12	0.48
1:A:72:PRO:HA	1:A:80:PHE:CE1	2.49	0.47
1:A:230:GLY:H	1:A:300:GLN:H	1.60	0.47
1:B:252:TYR:HB2	1:B:346:ILE:CG2	2.44	0.47
1:B:196:PRO:C	1:B:198:HIS:N	2.64	0.47
1:A:142:ILE:HG22	1:B:26:LEU:HA	1.96	0.47
1:B:274:THR:HG22	1:B:298:ASP:HA	1.95	0.47
1:B:193:LEU:HD21	1:B:198:HIS:HB3	1.96	0.47
1:A:251:ASP:CG	1:A:345:GLU:OE2	2.57	0.47
1:B:193:LEU:HA	1:B:264:TYR:HE2	1.79	0.47
1:B:319:ALA:O	1:B:322:ALA:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:PRO:HA	1:B:256:PRO:HG2	1.97	0.46
1:B:194:LEU:HD11	1:B:196:PRO:CB	2.45	0.46
1:B:229:THR:HA	1:B:300:GLN:O	2.16	0.46
1:B:188:ALA:HB1	1:B:264:TYR:CE2	2.50	0.46
1:B:319:ALA:O	1:B:322:ALA:HB3	2.15	0.45
1:A:141:SER:OG	1:B:27:GLU:OE2	2.20	0.45
1:B:25:VAL:HG23	3:B:502:HOH:O	2.17	0.45
1:B:172:CYS:SG	1:B:310:MET:HG3	2.57	0.45
1:B:262:GLU:HG2	1:B:263:GLU:OE2	2.16	0.45
1:A:264:TYR:O	1:A:269:ALA:HB3	2.16	0.45
1:B:311:PRO:O	1:B:314:GLU:N	2.48	0.45
1:B:317:VAL:CG1	1:B:318:LYS:H	2.11	0.45
1:B:252:TYR:CD1	1:B:292:VAL:HG13	2.51	0.44
1:B:253:GLU:C	1:B:255:ILE:N	2.61	0.44
1:A:251:ASP:HA	1:A:345:GLU:HB2	2.00	0.44
1:B:166:SER:HA	1:B:167:THR:CB	2.47	0.44
1:B:188:ALA:HB1	1:B:264:TYR:HD2	1.82	0.44
1:A:181:GLY:O	1:A:184:VAL:HG22	2.18	0.44
1:B:194:LEU:CD1	1:B:196:PRO:CB	2.95	0.44
1:B:23:PRO:HB3	1:B:273:GLY:CA	2.46	0.44
1:B:241:CYS:HB3	1:B:246:VAL:O	2.17	0.44
1:B:169:GLN:NE2	1:B:223:LYS:HE2	2.33	0.43
1:B:172:CYS:HB2	1:B:334:LEU:HD12	1.99	0.43
1:B:175:PHE:HA	1:B:329:THR:O	2.17	0.43
1:B:206:MET:HE3	1:B:301:THR:HB	2.01	0.43
1:B:246:VAL:HG11	1:B:334:LEU:HB3	2.00	0.43
1:A:179:PRO:HA	1:A:256:PRO:HB2	2.00	0.43
1:B:41:ASN:C	1:B:43:LEU:N	2.77	0.43
1:B:285:MET:HE3	1:B:288:MET:HB2	2.00	0.42
1:A:34:GLN:N	1:A:35:ALA:HB2	2.31	0.42
1:A:102:ALA:C	1:A:103:ILE:HD12	2.44	0.42
1:B:72:PRO:HA	1:B:80:PHE:CE1	2.54	0.42
1:A:106:LEU:O	1:A:141:SER:HA	2.20	0.42
1:B:83:ILE:CG2	1:B:206:MET:HE2	2.48	0.42
1:B:23:PRO:HG3	1:B:272:GLY:O	2.20	0.42
1:A:288:MET:HE2	1:A:292:VAL:HG12	2.01	0.42
1:A:107:GLY:HA2	1:A:142:ILE:O	2.20	0.42
1:B:193:LEU:HD23	1:B:193:LEU:HA	1.53	0.42
1:B:247:GLN:HG2	1:B:343:MET:HA	2.02	0.42
1:B:15:ALA:N	1:B:75:ASP:OD1	2.53	0.41
1:B:110:ILE:H	1:B:110:ILE:HG12	1.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:ILE:HG22	1:A:328:LEU:HD13	2.02	0.41
1:A:234:LEU:CD1	1:A:296:LEU:HB3	2.50	0.41
1:A:290:ARG:HG2	1:A:293:ARG:NH2	2.35	0.41
1:B:46:ASN:CB	1:B:50:ASP:HB2	2.51	0.41
1:A:226:THR:HG22	1:A:240:MET:HE3	2.02	0.40
1:B:13:HIS:N	1:B:75:ASP:O	2.54	0.40
1:A:76:ASN:HA	1:A:77:PRO:HD3	1.97	0.40
1:B:201:LEU:HD12	1:B:201:LEU:O	2.21	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	317/347 (91%)	295 (93%)	14 (4%)	8 (2%)	4	2
1	B	333/347 (96%)	300 (90%)	22 (7%)	11 (3%)	3	1
All	All	650/694 (94%)	595 (92%)	36 (6%)	19 (3%)	3	1

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	35	ALA
1	A	37	PHE
1	A	38	VAL
1	A	40	PRO
1	A	41	ASN
1	A	346	ILE
1	B	312	GLU
1	B	314	GLU
1	B	317	VAL

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Mol	Chain	Res	Type
1	B	318	LYS
1	A	46	ASN
1	B	196	PRO
1	B	197	GLU
1	B	30	LEU
1	B	42	LEU
1	B	14	GLY
1	B	38	VAL
1	B	23	PRO

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	208/269 (77%)	187 (90%)	21 (10%)	7	5
1	B	228/269 (85%)	195 (86%)	33 (14%)	3	1
All	All	436/538 (81%)	382 (88%)	54 (12%)	4	3

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ILE
1	A	30	LEU
1	A	32	SER
1	A	34	GLN
1	A	43	LEU
1	A	48	THR
1	A	54	VAL
1	A	103	ILE
1	A	123	GLU
1	A	142	ILE
1	A	153	VAL
1	A	156	ILE
1	A	169	GLN
1	A	184	VAL

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Mol	Chain	Res	Type
1	A	203	THR
1	A	211	ILE
1	A	250	VAL
1	A	284	LEU
1	A	328	LEU
1	A	334	LEU
1	A	343	MET
1	B	21	ILE
1	B	25	VAL
1	B	28	THR
1	B	30	LEU
1	B	34	GLN
1	B	38	VAL
1	B	48	THR
1	B	51	ASP
1	B	70	PHE
1	B	73	ILE
1	B	110	ILE
1	B	114	SER
1	B	184	VAL
1	B	193	LEU
1	B	194	LEU
1	B	197	GLU
1	B	222	VAL
1	B	238	SER
1	B	242	GLN
1	B	246	VAL
1	B	258	LEU
1	B	261	VAL
1	B	262	GLU
1	B	276	ARG
1	B	288	MET
1	B	295	LEU
1	B	296	LEU
1	B	312	GLU
1	B	314	GLU
1	B	317	VAL
1	B	335	VAL
1	B	344	VAL
1	B	347	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	169	GLN
1	B	315	ASN

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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

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	319/347 (91%)	0.38	18 (5%) 30 28	55, 81, 134, 178	0
1	B	335/347 (96%)	0.56	29 (8%) 16 15	52, 90, 144, 168	0
All	All	654/694 (94%)	0.47	47 (7%) 21 20	52, 85, 141, 178	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	ILE	4.9
1	B	37	PHE	4.3
1	B	31	HIS	4.0
1	B	201	LEU	4.0
1	B	38	VAL	3.9
1	B	14	GLY	3.8
1	A	185	LEU	3.6
1	A	29	ILE	3.3
1	B	35	ALA	3.2
1	B	49	ARG	3.0
1	B	18	GLY	3.0
1	A	38	VAL	2.9
1	A	39	ASP	2.9
1	B	24	LYS	2.8
1	B	45	GLY	2.7
1	B	29	ILE	2.7
1	B	22	SER	2.7
1	A	269	ALA	2.6
1	A	347	ARG	2.6
1	B	46	ASN	2.6
1	A	339	GLY	2.5
1	B	48	THR	2.5
1	B	25	VAL	2.5
1	B	328	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	41	ASN	2.4
1	B	39	ASP	2.4
1	A	47	GLU	2.4
1	B	167	THR	2.3
1	B	40	PRO	2.3
1	A	49	ARG	2.3
1	B	47	GLU	2.3
1	A	346	ILE	2.2
1	A	197	GLU	2.2
1	B	264	TYR	2.2
1	A	31	HIS	2.2
1	B	34	GLN	2.2
1	A	36	LYS	2.1
1	B	17	CYS	2.1
1	A	338	ARG	2.1
1	A	30	LEU	2.1
1	B	15	ALA	2.1
1	B	335	VAL	2.1
1	B	308	ALA	2.1
1	A	137	ALA	2.0
1	A	42	LEU	2.0
1	B	165	ASN	2.0
1	A	73	ILE	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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	B	401	1/1	-0.02	0.33	148,148,148,148	0
2	MG	A	401	1/1	0.49	0.17	106,106,106,106	0

6.5 Other polymers [i](#)

There are no such residues in this entry.