



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:58 AM UTC

PDB ID : 3A9Z / pdb_00003a9z
Title : Crystal structure of ras selenocysteine lyase in complex with selenopropionate
Authors : Omi, R.; Hirotsu, K.
Deposited on : 2009-11-09
Resolution : 1.55 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

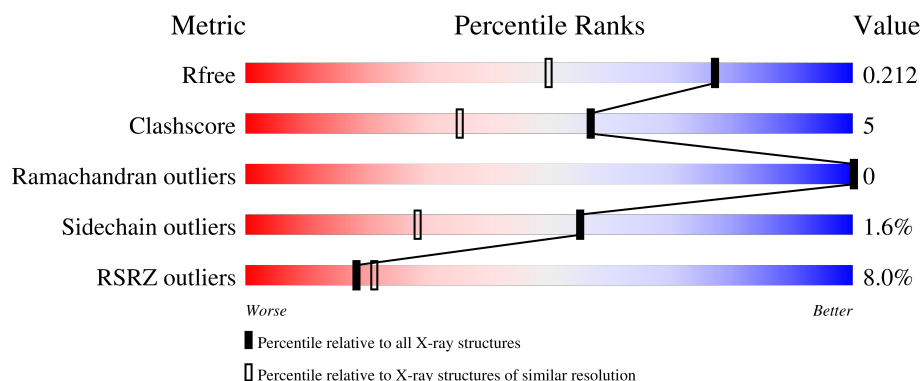
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	432	
1	B	432	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SLP	A	502[A]	-	X	-	-
3	SLP	A	502[B]	-	X	-	-
3	SLP	B	503[A]	-	X	-	-

2 Entry composition [i](#)

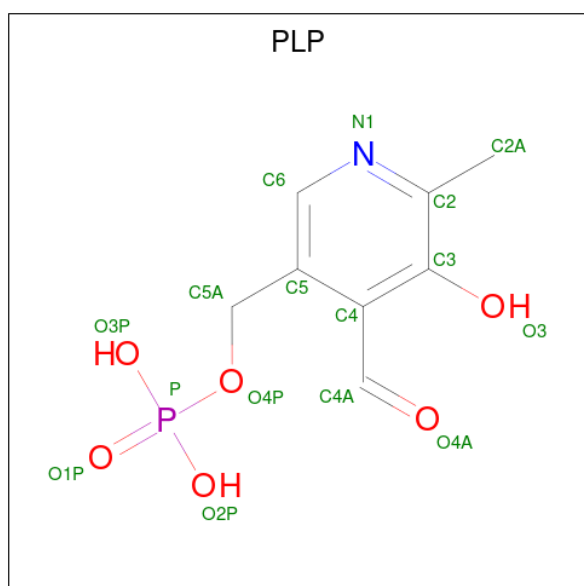
There are 5 unique types of molecules in this entry. The entry contains 6730 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Selenocysteine lyase.

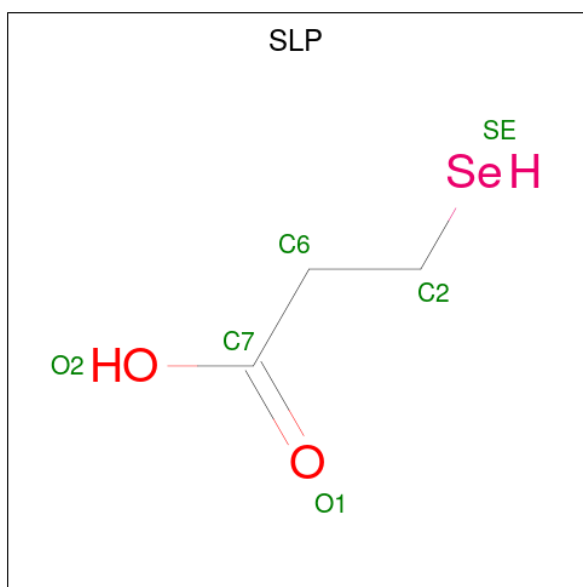
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	0	0
			3087	1929	556	585	17			
1	B	399	Total	C	N	O	S	0	0	0
			3052	1908	550	577	17			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



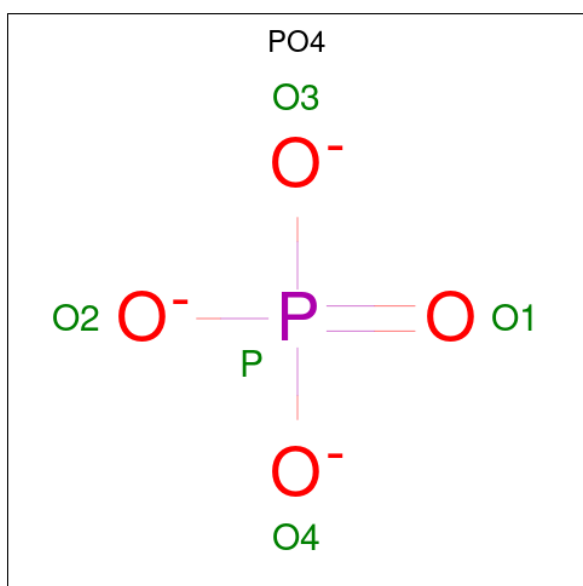
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is 3-selanylpropanoic acid (CCD ID: SLP) (formula: $C_3H_6O_2Se$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	Se	0	1
			12	6	4	2		
3	B	1	Total	C	O	Se	0	1
			12	6	4	2		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

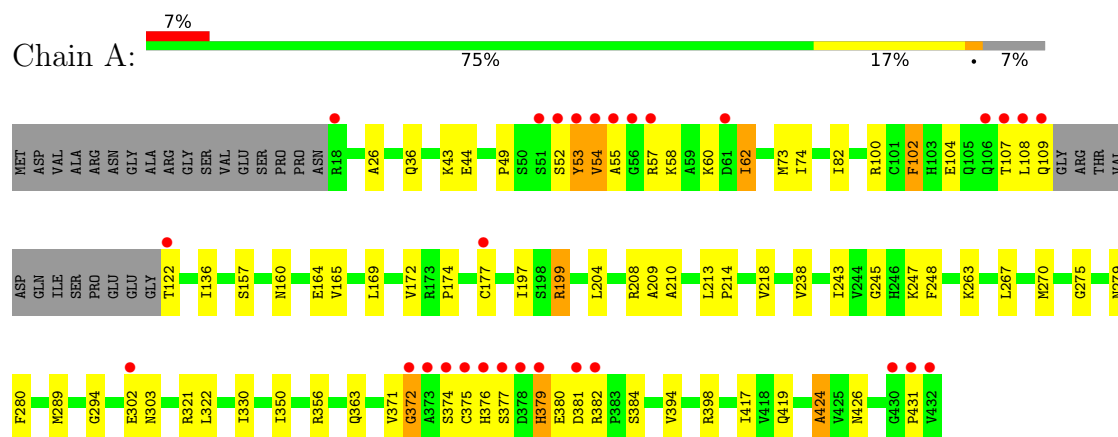
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	273	Total 273	O 273	0	0
5	B	254	Total 254	O 254	0	0

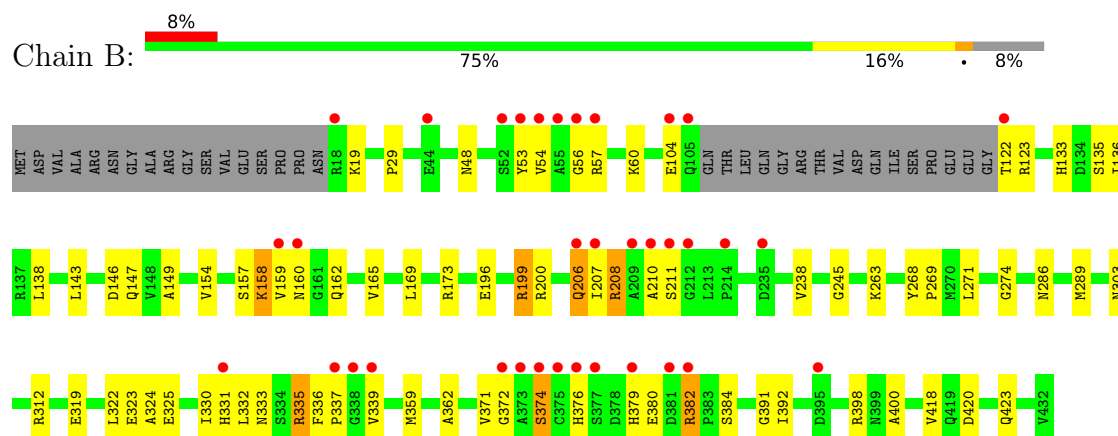
3 Residue-property plots [i](#)

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: Selenocysteine lyase



• Molecule 1: Selenocysteine lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.42Å 101.21Å 197.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 1.55 19.99 – 1.55	Depositor EDS
% Data completeness (in resolution range)	98.6 (19.99-1.55) 98.7 (19.99-1.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 1.55Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.187 , 0.209 0.189 , 0.212	Depositor DCC
R_{free} test set	15893 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.155	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6730	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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: SLP, PLP, 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	1.58	22/3139 (0.7%)	1.46	28/4259 (0.7%)
1	B	1.56	25/3104 (0.8%)	1.47	27/4213 (0.6%)
All	All	1.57	47/6243 (0.8%)	1.46	55/8472 (0.6%)

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	B	0	1

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	263	LYS	CA-CB	-8.70	1.38	1.54
1	A	431	PRO	C-O	8.15	1.33	1.23
1	A	73	MET	SD-CE	7.33	1.97	1.79
1	B	332	LEU	C-O	6.79	1.31	1.23
1	A	398	ARG	C-O	-6.76	1.15	1.24
1	A	53	TYR	C-O	6.76	1.31	1.23
1	B	208	ARG	C-O	6.70	1.32	1.24
1	A	177	CYS	CB-SG	-6.59	1.59	1.81
1	A	107	THR	C-O	6.58	1.31	1.24
1	B	359	MET	SD-CE	-6.37	1.63	1.79
1	A	350	ILE	CA-CB	-6.29	1.46	1.54
1	A	218	VAL	C-O	-6.21	1.17	1.24
1	A	424	ALA	CA-CB	6.19	1.62	1.53
1	B	104	GLU	C-O	6.16	1.32	1.24
1	A	417	ILE	CA-CB	5.93	1.61	1.54
1	A	43	LYS	CA-C	-5.84	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	243	ILE	C-O	-5.83	1.17	1.24
1	B	149	ALA	CA-CB	5.82	1.64	1.53
1	B	53	TYR	C-O	5.81	1.31	1.23
1	A	376	HIS	C-O	5.80	1.31	1.24
1	B	362	ALA	CA-CB	-5.74	1.44	1.53
1	B	382	ARG	C-O	5.67	1.31	1.24
1	A	294	GLY	C-O	-5.61	1.17	1.23
1	B	206	GLN	C-O	5.57	1.30	1.24
1	B	312	ARG	CZ-NH1	5.53	1.40	1.32
1	B	376	HIS	C-O	5.50	1.31	1.24
1	B	324	ALA	CA-CB	-5.48	1.44	1.53
1	B	286	ASN	C-O	-5.46	1.16	1.23
1	A	426	ASN	C-O	5.45	1.30	1.24
1	B	206	GLN	CG-CD	5.44	1.65	1.52
1	A	379	HIS	C-O	5.36	1.29	1.23
1	B	271	LEU	C-O	-5.29	1.17	1.23
1	B	210	ALA	C-O	5.21	1.30	1.24
1	B	374	SER	C-O	5.16	1.30	1.24
1	B	173	ARG	N-CA	5.12	1.49	1.46
1	B	263	LYS	CA-C	5.11	1.57	1.52
1	A	54	VAL	CA-C	5.11	1.59	1.52
1	A	248	PHE	CA-CB	5.10	1.60	1.53
1	B	56	GLY	C-O	5.09	1.30	1.23
1	A	62	ILE	CA-CB	-5.08	1.48	1.54
1	B	339	VAL	CA-C	-5.08	1.46	1.52
1	A	58	LYS	N-CA	5.08	1.52	1.46
1	A	214	PRO	CA-C	-5.06	1.45	1.52
1	B	159	VAL	CA-CB	5.03	1.60	1.55
1	B	162	GLN	CA-C	-5.03	1.46	1.52
1	A	36	GLN	CD-OE1	5.02	1.33	1.23
1	B	57	ARG	C-O	5.01	1.30	1.24

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	136	ILE	N-CA-C	-9.72	103.94	111.62
1	B	19	LYS	N-CA-C	-9.00	97.54	110.59
1	B	154	VAL	CA-C-N	-8.36	111.40	119.85
1	B	154	VAL	C-N-CA	-8.36	111.40	119.85
1	A	52	SER	CB-CA-C	-7.28	101.61	111.88
1	A	384	SER	N-CA-C	6.85	119.44	109.71
1	B	54	VAL	N-CA-C	6.73	118.29	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	245	GLY	N-CA-C	6.62	120.91	112.83
1	B	238	VAL	N-CA-C	6.60	117.24	110.05
1	A	245	GLY	N-CA-C	6.57	120.84	112.83
1	A	136	ILE	N-CA-C	-6.55	105.92	111.56
1	A	371	VAL	CA-C-N	-6.51	108.64	121.41
1	A	371	VAL	C-N-CA	-6.51	108.64	121.41
1	B	54	VAL	N-CA-CB	6.26	118.30	110.47
1	A	74	ILE	N-CA-C	6.12	117.98	112.17
1	A	321	ARG	CG-CD-NE	-6.10	98.57	112.00
1	A	53	TYR	N-CA-C	-6.05	98.94	108.14
1	A	394	VAL	N-CA-C	5.95	117.40	110.62
1	A	199	ARG	CB-CG-CD	-5.83	97.90	111.30
1	A	210	ALA	N-CA-C	-5.69	105.08	111.28
1	B	54	VAL	CB-CA-C	-5.62	104.68	112.04
1	B	29	PRO	N-CA-C	-5.61	103.36	111.22
1	A	102	PHE	N-CA-C	-5.61	105.07	111.07
1	A	58	LYS	CG-CD-CE	5.55	124.07	111.30
1	B	48	ASN	CA-C-N	-5.51	114.42	119.82
1	B	48	ASN	C-N-CA	-5.51	114.42	119.82
1	B	274	GLY	N-CA-C	5.50	126.21	113.18
1	B	400	ALA	N-CA-C	5.45	118.35	110.28
1	B	384	SER	CA-C-N	-5.43	114.07	119.56
1	B	384	SER	C-N-CA	-5.43	114.07	119.56
1	B	372	GLY	N-CA-C	-5.42	103.24	111.42
1	A	381	ASP	CA-C-N	-5.41	114.09	121.61
1	A	381	ASP	C-N-CA	-5.41	114.09	121.61
1	A	372	GLY	N-CA-C	-5.38	100.42	113.18
1	A	107	THR	N-CA-C	-5.37	105.11	110.97
1	A	109	GLN	N-CA-C	-5.36	95.99	111.00
1	B	335	ARG	CB-CA-C	-5.36	104.40	111.74
1	A	238	VAL	N-CA-C	5.31	115.83	110.05
1	A	164	GLU	CA-CB-CG	-5.27	103.56	114.10
1	A	43	LYS	N-CA-C	5.26	118.16	111.69
1	A	270	MET	N-CA-C	-5.26	106.01	112.90
1	B	303	ASN	N-CA-C	5.22	120.80	113.02
1	B	269	PRO	N-CA-C	5.17	119.60	111.38
1	A	44	GLU	N-CA-C	5.15	119.83	113.50
1	B	392	ILE	N-CA-C	-5.15	104.20	108.63
1	B	138	LEU	CA-C-N	-5.11	113.77	119.19
1	B	138	LEU	C-N-CA	-5.11	113.77	119.19
1	B	199	ARG	NE-CZ-NH1	-5.10	116.40	121.50
1	A	247	LYS	N-CA-C	-5.08	107.13	113.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	398	ARG	NE-CZ-NH1	5.06	126.56	121.50
1	A	197	ILE	N-CA-C	-5.06	105.56	110.42
1	A	209	ALA	O-C-N	5.05	127.47	122.12
1	B	123	ARG	CA-C-N	-5.03	114.77	119.85
1	B	123	ARG	C-N-CA	-5.03	114.77	119.85
1	A	164	GLU	N-CA-C	5.00	117.97	109.76

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	268	TYR	Sidechain

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	3087	0	3096	35	0
1	B	3052	0	3060	22	0
2	A	15	0	6	1	0
2	B	15	0	6	1	0
3	A	12	0	4	2	0
3	B	12	0	4	1	0
4	A	10	0	0	0	0
5	A	273	0	0	4	0
5	B	254	0	0	0	0
All	All	6730	0	6176	57	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 5.

All (57) 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)
2:A:501:PLP:C4A	3:A:502[A]:SLP:SE	2.28	1.30

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:501:PLP:C4A	3:B:503[A]:SLP:SE	2.50	1.09
1:A:379:HIS:HB3	1:A:382:ARG:HD2	1.52	0.88
1:A:102:PHE:HE2	1:A:122:THR:O	1.57	0.86
1:B:200:ARG:HG3	1:B:200:ARG:NH1	1.98	0.78
1:B:319:GLU:O	1:B:323:GLU:HG3	1.85	0.77
1:A:82:ILE:HD13	1:A:267:LEU:CD2	2.19	0.72
1:A:82:ILE:CD1	1:A:267:LEU:HD22	2.22	0.70
1:A:165:VAL:O	1:A:169:LEU:HD23	1.91	0.70
1:A:379:HIS:CD2	1:A:382:ARG:CZ	2.75	0.70
1:B:200:ARG:HG3	1:B:200:ARG:HH11	1.57	0.69
1:A:377:SER:O	1:A:380:GLU:HG3	1.95	0.67
1:B:200:ARG:HH11	1:B:200:ARG:CG	2.09	0.65
1:A:104:GLU:O	1:A:108:LEU:HG	2.00	0.61
1:A:372:GLY:HA3	5:A:667:HOH:O	2.00	0.60
1:B:157:SER:HB3	1:B:160:ASN:OD1	2.02	0.59
1:B:165:VAL:O	1:B:169:LEU:HD23	2.02	0.59
1:B:325:GLU:HG3	1:B:418:VAL:CG1	2.34	0.58
1:B:325:GLU:HG3	1:B:418:VAL:HG11	1.86	0.57
1:A:54:VAL:HG23	1:A:57:ARG:NH2	2.19	0.57
1:A:82:ILE:CD1	1:A:267:LEU:CD2	2.81	0.56
1:A:267:LEU:HD23	1:A:280:PHE:HB3	1.88	0.55
1:A:54:VAL:HG23	1:A:57:ARG:HH22	1.72	0.54
1:B:133:HIS:HD1	1:B:135:SER:H	1.55	0.54
1:B:379:HIS:ND1	1:B:382:ARG:CZ	2.70	0.54
1:A:49:PRO:O	1:A:60:LYS:HE2	2.08	0.54
1:A:356:ARG:HD3	5:A:656:HOH:O	2.08	0.53
1:A:53:TYR:CE1	1:A:55:ALA:HB3	2.44	0.53
1:A:379:HIS:HD2	1:A:382:ARG:CZ	2.21	0.52
1:B:420:ASP:O	1:B:423:GLN:HG3	2.10	0.52
1:A:100:ARG:NH1	1:B:146:ASP:OD2	2.33	0.51
1:A:26:ALA:HB1	1:A:374:SER:OG	2.11	0.51
1:A:172:VAL:HG11	1:A:204:LEU:HD21	1.94	0.50
1:A:275:GLY:HA2	1:A:279:ASN:CG	2.39	0.48
1:A:82:ILE:HD13	1:A:267:LEU:HD22	1.90	0.48
1:B:379:HIS:CE1	1:B:382:ARG:NH2	2.83	0.46
1:A:322:LEU:HB3	1:A:330:ILE:HD13	1.96	0.46
1:B:196:GLU:OE2	1:B:199:ARG:NH2	2.44	0.46
1:A:289:MET:HE3	1:B:289:MET:HE3	1.97	0.46
1:B:158:LYS:HD3	1:B:391:GLY:HA2	1.99	0.45
1:B:207:ILE:O	1:B:208:ARG:C	2.59	0.44
1:B:371:VAL:HG21	1:B:380:GLU:HB3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:SER:HB3	1:A:160:ASN:OD1	2.18	0.44
1:A:302:GLU:HG3	1:A:303:ASN:CG	2.42	0.44
1:B:322:LEU:HB3	1:B:330:ILE:HD13	1.99	0.44
1:A:208:ARG:HB3	1:A:213:LEU:HB2	2.00	0.42
1:A:199:ARG:HG2	5:A:524:HOH:O	2.19	0.42
1:B:331:HIS:ND1	1:B:333:ASN:OD1	2.46	0.42
1:B:336:PHE:HB3	1:B:337:PRO:HD2	2.02	0.42
1:A:53:TYR:CD1	1:A:55:ALA:HB3	2.55	0.42
1:A:419:GLN:HG2	5:A:663:HOH:O	2.20	0.41
1:A:375:CYS:SG	3:A:502[B]:SLP:SE	3.29	0.41
1:A:275:GLY:HA2	1:A:279:ASN:ND2	2.36	0.41
1:A:53:TYR:HE1	1:A:55:ALA:HB3	1.87	0.41
1:A:102:PHE:CE2	1:A:122:THR:O	2.50	0.40
1:A:363:GLN:HB2	1:A:424:ALA:HB1	2.03	0.40
1:B:143:LEU:HD23	1:B:143:LEU:HA	1.89	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	399/432 (92%)	388 (97%)	11 (3%)	0	100	100
1	B	395/432 (91%)	384 (97%)	11 (3%)	0	100	100
All	All	794/864 (92%)	772 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	337/364 (93%)	334 (99%)	3 (1%)	70	51
1	B	333/364 (92%)	325 (98%)	8 (2%)	43	15
All	All	670/728 (92%)	659 (98%)	11 (2%)	55	28

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	ILE
1	A	174	PRO
1	A	263	LYS
1	B	60	LYS
1	B	122	THR
1	B	147	GLN
1	B	158	LYS
1	B	206	GLN
1	B	211	SER
1	B	335	ARG
1	B	374	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	279	ASN
1	A	363	GLN
1	A	365	GLN
1	A	376	HIS
1	A	379	HIS
1	B	303	ASN
1	B	354	GLN

5.3.3 RNA ⓘ

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](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	SLP	B	503[B]	-	3,5,5	2.82	2 (66%)	5,5,5	1.55	1 (20%)
3	SLP	A	502[B]	-	3,5,5	2.80	3 (100%)	5,5,5	1.54	1 (20%)
4	PO4	A	504	-	4,4,4	1.87	2 (50%)	6,6,6	0.82	0
2	PLP	A	501	1	15,15,16	1.84	3 (20%)	21,22,23	1.55	5 (23%)
3	SLP	B	503[A]	-	3,5,5	2.67	2 (66%)	5,5,5	3.31	3 (60%)
3	SLP	A	502[A]	-	3,5,5	2.53	2 (66%)	5,5,5	3.05	2 (40%)
4	PO4	A	505	-	4,4,4	1.83	1 (25%)	6,6,6	0.63	0
2	PLP	B	501	1	15,15,16	1.39	3 (20%)	21,22,23	1.11	2 (9%)

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	SLP	B	503[B]	-	-	2/3/3/3	-
3	SLP	A	502[B]	-	-	2/3/3/3	-
2	PLP	A	501	1	-	0/6/6/8	0/1/1/1
3	SLP	B	503[A]	-	-	2/3/3/3	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SLP	A	502[A]	-	-	2/3/3/3	-
2	PLP	B	501	1	-	0/6/6/8	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	PLP	P-O1P	5.39	1.67	1.50
3	B	503[A]	SLP	O2-C7	-3.76	1.18	1.30
3	B	503[B]	SLP	O2-C7	-3.71	1.18	1.30
3	A	502[B]	SLP	O1-C7	3.33	1.33	1.22
3	A	502[A]	SLP	O1-C7	3.21	1.32	1.22
2	B	501	PLP	C3-C2	-3.16	1.37	1.41
4	A	505	PO4	P-O3	-2.64	1.46	1.54
3	A	502[B]	SLP	C6-C7	2.62	1.56	1.50
4	A	504	PO4	P-O3	-2.56	1.47	1.54
2	A	501	PLP	C2A-C2	2.54	1.54	1.50
3	B	503[B]	SLP	O1-C7	2.53	1.30	1.22
2	B	501	PLP	C2A-C2	2.50	1.54	1.50
3	A	502[B]	SLP	O2-C7	-2.36	1.23	1.30
4	A	504	PO4	P-O2	-2.36	1.47	1.54
3	B	503[A]	SLP	O1-C7	2.33	1.29	1.22
3	A	502[A]	SLP	O2-C7	-2.29	1.23	1.30
2	A	501	PLP	P-O4P	-2.18	1.53	1.60
2	B	501	PLP	P-O1P	2.01	1.56	1.50

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	503[A]	SLP	SE-C2-C6	-6.20	97.54	112.77
3	A	502[A]	SLP	SE-C2-C6	-6.12	97.72	112.77
2	A	501	PLP	O4P-C5A-C5	4.09	117.03	109.36
3	B	503[A]	SLP	O2-C7-C6	2.60	122.22	114.00
2	A	501	PLP	C4A-C4-C5	-2.60	118.26	120.94
3	B	503[B]	SLP	O2-C7-C6	2.55	122.04	114.00
2	B	501	PLP	O4P-C5A-C5	2.37	113.80	109.36
2	B	501	PLP	C5-C6-N1	-2.29	120.10	123.83
3	A	502[A]	SLP	O2-C7-C6	2.21	121.00	114.00
2	A	501	PLP	C6-C5-C4	2.20	119.90	118.10
3	B	503[A]	SLP	C2-C6-C7	-2.17	108.22	113.28
2	A	501	PLP	C5-C6-N1	-2.14	120.34	123.83
3	A	502[B]	SLP	O2-C7-C6	2.10	120.63	114.00
2	A	501	PLP	C2A-C2-C3	2.07	123.22	120.80

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	503[A]	SLP	C2-C6-C7-O1
3	B	503[B]	SLP	C2-C6-C7-O1
3	A	502[B]	SLP	C2-C6-C7-O1
3	B	503[A]	SLP	C2-C6-C7-O2
3	A	502[B]	SLP	C2-C6-C7-O2
3	B	503[B]	SLP	C2-C6-C7-O2
3	A	502[A]	SLP	C2-C6-C7-O2
3	A	502[A]	SLP	C2-C6-C7-O1

There are no ring outliers.

5 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502[B]	SLP	1	0
2	A	501	PLP	1	0
3	B	503[A]	SLP	1	0
3	A	502[A]	SLP	1	0
2	B	501	PLP	1	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 ⓘ

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	403/432 (93%)	0.24	29 (7%) 21 25	13, 19, 36, 48	0
1	B	399/432 (92%)	0.42	35 (8%) 15 17	14, 22, 37, 46	0
All	All	802/864 (92%)	0.33	64 (7%) 18 21	13, 20, 37, 48	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	432	VAL	8.4
1	A	373	ALA	8.0
1	A	56	GLY	6.8
1	B	54	VAL	5.6
1	A	122	THR	5.4
1	A	379	HIS	5.3
1	B	373	ALA	5.3
1	A	431	PRO	5.2
1	A	374	SER	5.1
1	A	375	CYS	5.0
1	B	56	GLY	4.7
1	A	108	LEU	4.5
1	A	381	ASP	4.5
1	B	375	CYS	4.4
1	A	52	SER	4.4
1	A	377	SER	4.3
1	B	379	HIS	4.0
1	B	210	ALA	4.0
1	B	105	GLN	3.9
1	B	209	ALA	3.8
1	B	377	SER	3.7
1	B	376	HIS	3.7
1	A	376	HIS	3.7
1	A	18	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	53	TYR	3.6
1	B	337	PRO	3.6
1	A	372	GLY	3.5
1	B	372	GLY	3.4
1	B	382	ARG	3.4
1	A	57	ARG	3.3
1	A	378	ASP	3.2
1	B	207	ILE	3.1
1	B	159	VAL	3.0
1	A	54	VAL	2.9
1	B	57	ARG	2.9
1	A	55	ALA	2.9
1	B	122	THR	2.9
1	B	331	HIS	2.8
1	A	106	GLN	2.8
1	A	382	ARG	2.7
1	B	55	ALA	2.7
1	B	374	SER	2.7
1	A	61	ASP	2.7
1	B	395	ASP	2.7
1	B	18	ARG	2.6
1	B	235	ASP	2.6
1	B	104	GLU	2.6
1	B	338	GLY	2.6
1	B	44	GLU	2.5
1	A	109	GLN	2.5
1	B	206	GLN	2.5
1	B	214	PRO	2.4
1	B	53	TYR	2.3
1	B	212	GLY	2.3
1	A	302	GLU	2.3
1	A	51	SER	2.3
1	A	430	GLY	2.1
1	B	52	SER	2.1
1	B	339	VAL	2.1
1	B	211	SER	2.1
1	A	107	THR	2.1
1	B	381	ASP	2.1
1	B	160	ASN	2.1
1	A	177	CYS	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
4	PO4	A	505	5/5	0.89	0.11	36,38,40,40	0
4	PO4	A	504	5/5	0.94	0.10	29,29,30,31	0
3	SLP	B	503[A]	6/6	0.96	0.15	19,23,24,25	6
3	SLP	B	503[B]	6/6	0.96	0.15	9,18,19,20	6
2	PLP	A	501	15/16	0.98	0.05	13,15,17,18	0
2	PLP	B	501	15/16	0.98	0.05	16,18,19,21	0
3	SLP	A	502[A]	6/6	0.98	0.13	13,20,21,21	6
3	SLP	A	502[B]	6/6	0.98	0.13	10,18,19,20	6

6.5 Other polymers [i](#)

There are no such residues in this entry.