



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:20 PM EDT

PDB ID : 7F1P / pdb_00007f1p
Title : Crystal structure of Pseudomonas putida methionine gamma-lyase Q349S mutant ligand-free form.
Authors : Okawa, A.; Handa, H.; Yasuda, E.; Murota, M.; Kudo, D.; Tamura, T.; Shiba, T.; Inagaki, K.
Deposited on : 2021-06-09
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

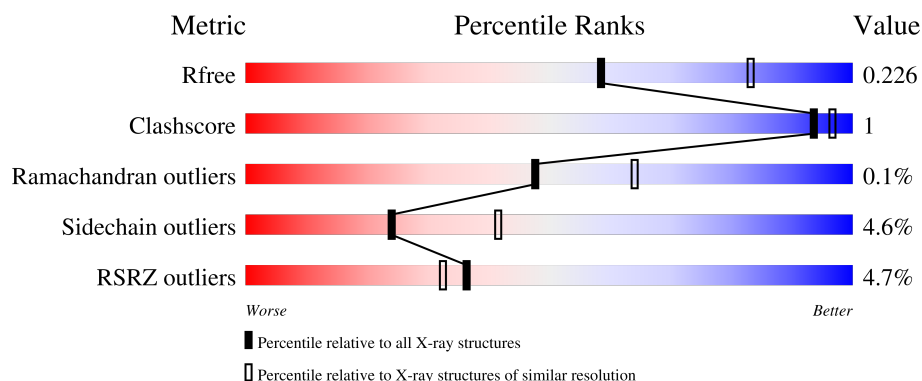
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

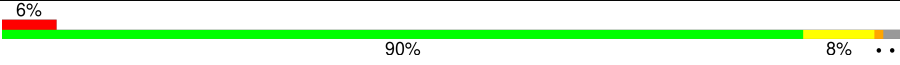
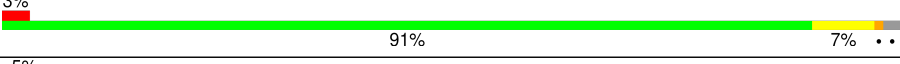
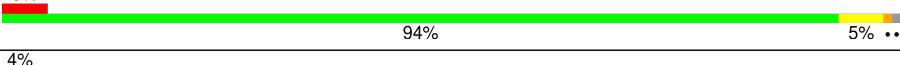
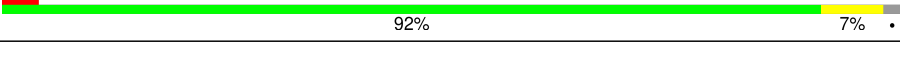
The reported resolution of this entry is 2.40 Å.

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))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (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	398	 6% 90% 8% ..
1	B	398	 3% 91% 7% ..
1	C	398	 5% 94% 5% ..
1	D	398	 4% 92% 7% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12003 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 L-methionine gamma-lyase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	P	S	0	0	0
			2963	1869	522	554	1	17			
1	B	392	Total	C	N	O	P	S	0	0	0
			2963	1869	522	554	1	17			
1	C	396	Total	C	N	O	P	S	0	0	0
			2990	1884	528	560	1	17			
1	D	392	Total	C	N	O	P	S	0	0	0
			2963	1869	522	554	1	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	349	SER	GLN	engineered mutation	UNP P13254
B	349	SER	GLN	engineered mutation	UNP P13254
C	349	SER	GLN	engineered mutation	UNP P13254
D	349	SER	GLN	engineered mutation	UNP P13254

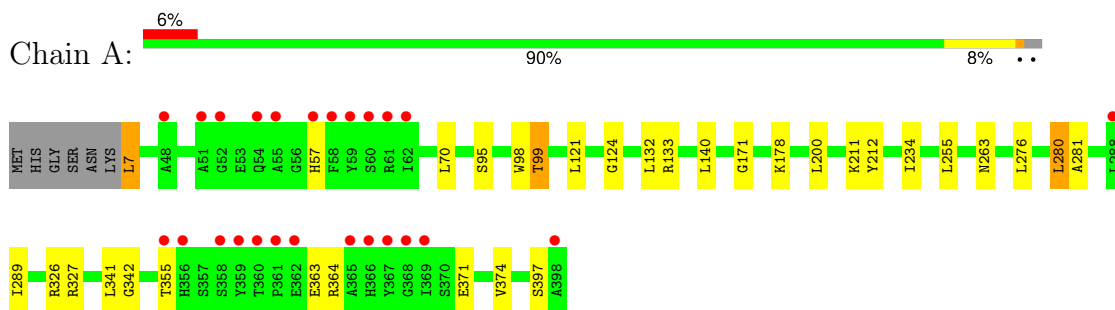
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	17	Total	O	0	0
			17	17		
2	B	43	Total	O	0	0
			43	43		
2	C	40	Total	O	0	0
			40	40		
2	D	24	Total	O	0	0
			24	24		

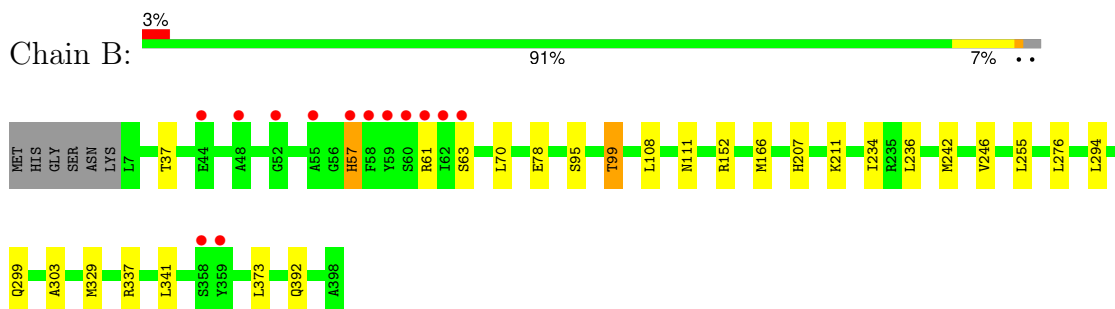
3 Residue-property plots [i](#)

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.

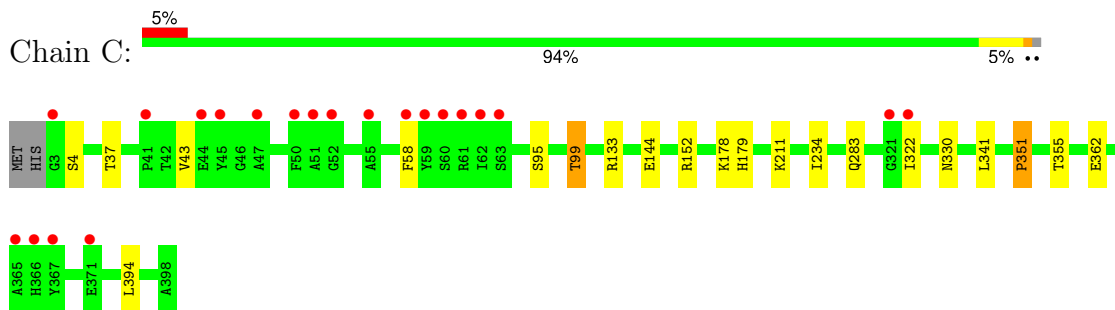
- Molecule 1: L-methionine gamma-lyase



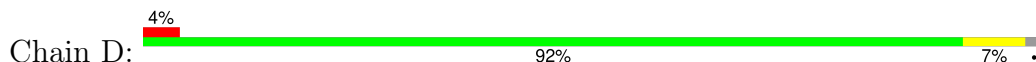
- Molecule 1: L-methionine gamma-lyase

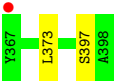
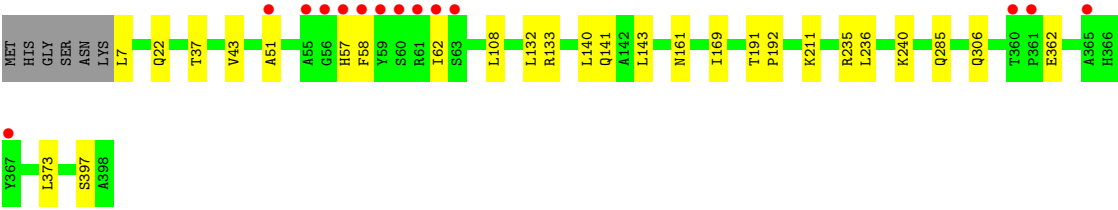


- Molecule 1: L-methionine gamma-lyase



- Molecule 1: L-methionine gamma-lyase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	154.14Å 153.01Å 80.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.95 – 2.40 19.95 – 2.40	Depositor EDS
% Data completeness (in resolution range)	79.9 (19.95-2.40) 79.9 (19.95-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.184 , 0.227 0.186 , 0.226	Depositor DCC
R_{free} test set	3085 reflections (4.10%)	wwPDB-VP
Wilson B-factor (Å ²)	38.9	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 18.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12003	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% 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: LLP

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	0.66	0/3002	0.86	0/4075
1	B	0.70	0/3002	0.91	1/4075 (0.0%)
1	C	0.69	0/3029	0.87	2/4110 (0.0%)
1	D	0.66	0/3002	0.89	0/4075
All	All	0.68	0/12035	0.88	3/16335 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	HIS	N-CA-C	7.14	118.02	108.38
1	C	283	GLN	CA-C-N	5.66	125.13	119.24
1	C	283	GLN	C-N-CA	5.66	125.13	119.24

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	2963	0	2922	13	0
1	B	2963	0	2922	11	0
1	C	2990	0	2949	7	0
1	D	2963	0	2922	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	17	0	0	1	0
2	B	43	0	0	0	0
2	C	40	0	0	1	0
2	D	24	0	0	0	0
All	All	12003	0	11715	33	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 1.

All (33) 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:99:THR:HG21	1:A:234:ILE:HA	1.72	0.71
1:C:99:THR:HG21	1:C:234:ILE:HA	1.73	0.70
1:B:166:MET:HE1	1:B:303:ALA:HA	1.75	0.69
1:A:95:SER:O	1:A:99:THR:HG23	1.99	0.62
1:B:99:THR:HG21	1:B:234:ILE:HA	1.81	0.61
1:D:62:ILE:O	1:D:235:ARG:NH1	2.36	0.58
1:D:169:ILE:H	1:D:306:GLN:HE22	1.51	0.56
1:B:61:ARG:HG3	1:B:246:VAL:HG21	1.88	0.56
1:B:70:LEU:HD21	1:B:255:LEU:HD23	1.90	0.54
1:C:95:SER:O	1:C:99:THR:CG2	2.55	0.54
1:A:98:TRP:HZ2	1:B:242:MET:HE2	1.73	0.52
1:A:140:LEU:HD11	1:A:171:GLY:HA3	1.91	0.52
1:C:330:ASN:HD22	1:C:330:ASN:N	2.08	0.51
1:B:166:MET:H	1:B:299:GLN:HE22	1.58	0.51
1:B:166:MET:HE1	1:B:303:ALA:CA	2.44	0.46
1:C:144:GLU:OE2	1:C:179:HIS:NE2	2.46	0.45
1:D:236:LEU:O	1:D:240:LYS:HE3	2.17	0.45
1:C:95:SER:O	1:C:99:THR:HG23	2.17	0.44
1:A:7:LEU:N	2:A:401:HOH:O	2.50	0.43
1:B:329:MET:HE3	1:B:337:ARG:HG2	2.00	0.43
1:A:70:LEU:HD21	1:A:255:LEU:HD23	2.00	0.43
1:A:281:ALA:HA	1:A:289:ILE:HD11	2.00	0.43
1:A:276:LEU:O	1:A:280:LEU:HD22	2.19	0.42
1:A:124:GLY:CA	1:B:242:MET:HE3	2.50	0.42
1:A:327:ARG:NH2	1:A:397:SER:O	2.53	0.42
1:D:191:THR:HB	1:D:192:PRO:HD2	2.03	0.41
1:A:121:LEU:C	1:A:132:LEU:HD21	2.46	0.41
1:C:330:ASN:HD21	1:D:43:VAL:HG22	1.86	0.41
1:B:95:SER:O	1:B:99:THR:CG2	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:TYR:CE1	1:A:342:GLY:HA2	2.56	0.41
1:C:355:THR:HG23	2:C:440:HOH:O	2.21	0.41
1:B:78:GLU:OE2	1:B:207:HIS:CE1	2.74	0.40
1:A:364:ARG:NH2	1:A:371:GLU:OE2	2.50	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	389/398 (98%)	378 (97%)	11 (3%)	0	100	100
1	B	389/398 (98%)	379 (97%)	10 (3%)	0	100	100
1	C	393/398 (99%)	382 (97%)	10 (2%)	1 (0%)	36	50
1	D	389/398 (98%)	372 (96%)	16 (4%)	1 (0%)	36	50
All	All	1560/1592 (98%)	1511 (97%)	47 (3%)	2 (0%)	48	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	51	ALA
1	C	351	PRO

5.3.2 Protein sidechains [i](#)

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	301/306 (98%)	288 (96%)	13 (4%)	26	44
1	B	301/306 (98%)	288 (96%)	13 (4%)	26	44
1	C	304/306 (99%)	291 (96%)	13 (4%)	26	44
1	D	301/306 (98%)	285 (95%)	16 (5%)	20	36
All	All	1207/1224 (99%)	1152 (95%)	55 (5%)	24	41

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	57	HIS
1	A	99	THR
1	A	133	ARG
1	A	178	LYS
1	A	200	LEU
1	A	263	ASN
1	A	280	LEU
1	A	326	ARG
1	A	341	LEU
1	A	355	THR
1	A	363	GLU
1	A	374	VAL
1	B	37	THR
1	B	57	HIS
1	B	63	SER
1	B	99	THR
1	B	108	LEU
1	B	111	ASN
1	B	152	ARG
1	B	236	LEU
1	B	276	LEU
1	B	294	LEU
1	B	341	LEU
1	B	373	LEU
1	B	392	GLN
1	C	4	SER
1	C	37	THR
1	C	43	VAL
1	C	58	PHE
1	C	99	THR
1	C	133	ARG
1	C	152	ARG

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Mol	Chain	Res	Type
1	C	178	LYS
1	C	322	ILE
1	C	341	LEU
1	C	351	PRO
1	C	362	GLU
1	C	394	LEU
1	D	7	LEU
1	D	22	GLN
1	D	37	THR
1	D	57	HIS
1	D	58	PHE
1	D	108	LEU
1	D	132	LEU
1	D	133	ARG
1	D	140	LEU
1	D	141	GLN
1	D	143	LEU
1	D	161	ASN
1	D	285	GLN
1	D	362	GLU
1	D	373	LEU
1	D	397	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	161	ASN
1	A	250	HIS
1	A	263	ASN
1	A	283	GLN
1	A	333	GLN
1	A	391	GLN
1	B	68	ASN
1	B	195	GLN
1	B	228	GLN
1	B	290	HIS
1	B	299	GLN
1	B	306	GLN
1	B	330	ASN
1	B	333	GLN
1	B	366	HIS
1	B	391	GLN

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Mol	Chain	Res	Type
1	B	392	GLN
1	C	187	ASN
1	C	207	HIS
1	C	228	GLN
1	C	330	ASN
1	D	68	ASN
1	D	161	ASN
1	D	187	ASN
1	D	299	GLN
1	D	306	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	LLP	B	211	1	23,24,25	2.42	5 (21%)	25,32,34	1.44	6 (24%)
1	LLP	D	211	1	23,24,25	2.43	5 (21%)	25,32,34	1.46	7 (28%)
1	LLP	A	211	1	23,24,25	2.50	5 (21%)	25,32,34	1.41	6 (24%)
1	LLP	C	211	1	23,24,25	2.37	5 (21%)	25,32,34	1.42	5 (20%)

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
1	LLP	B	211	1	-	4/16/17/19	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	LLP	D	211	1	-	4/16/17/19	0/1/1/1
1	LLP	A	211	1	-	4/16/17/19	0/1/1/1
1	LLP	C	211	1	-	4/16/17/19	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	211	LLP	C3-C2	7.06	1.48	1.41
1	B	211	LLP	C3-C2	6.88	1.48	1.41
1	D	211	LLP	C3-C2	6.59	1.47	1.41
1	C	211	LLP	C3-C2	6.44	1.47	1.41
1	D	211	LLP	C4'-NZ	5.63	1.46	1.27
1	C	211	LLP	C4'-NZ	5.58	1.45	1.27
1	A	211	LLP	C4'-NZ	5.56	1.45	1.27
1	B	211	LLP	C4'-NZ	5.51	1.45	1.27
1	B	211	LLP	C4-C5	5.39	1.49	1.42
1	A	211	LLP	C4-C5	5.19	1.49	1.42
1	D	211	LLP	C4-C5	5.16	1.49	1.42
1	C	211	LLP	C4-C5	4.94	1.48	1.42
1	A	211	LLP	C4-C3	4.71	1.48	1.41
1	D	211	LLP	C4-C3	4.47	1.48	1.41
1	C	211	LLP	C4-C3	4.30	1.48	1.41
1	B	211	LLP	C4-C3	3.92	1.47	1.41
1	B	211	LLP	C4-C4'	2.75	1.52	1.46
1	D	211	LLP	C4-C4'	2.56	1.52	1.46
1	C	211	LLP	C4-C4'	2.46	1.51	1.46
1	A	211	LLP	C4-C4'	2.20	1.51	1.46

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	211	LLP	C4-C3-C2	-3.05	118.42	120.14
1	B	211	LLP	C6-N1-C2	2.66	124.03	119.20
1	C	211	LLP	OP4-C5'-C5	2.63	114.29	109.36
1	D	211	LLP	C4-C3-C2	-2.62	118.67	120.14
1	C	211	LLP	C3-C4-C5	-2.53	116.24	118.28
1	A	211	LLP	C4-C3-C2	-2.48	118.75	120.14
1	D	211	LLP	C6-N1-C2	2.47	123.68	119.20
1	D	211	LLP	OP4-P-OP1	-2.45	99.81	106.44
1	A	211	LLP	C6-N1-C2	2.39	123.53	119.20
1	A	211	LLP	OP3-P-OP2	2.39	116.76	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	211	LLP	C4-C4'-NZ	-2.34	113.24	124.04
1	A	211	LLP	C4-C4'-NZ	-2.33	113.27	124.04
1	A	211	LLP	OP3-P-OP4	-2.31	100.64	106.67
1	C	211	LLP	OP3-P-OP2	2.29	116.38	107.80
1	D	211	LLP	CD-CE-NZ	2.27	116.85	110.83
1	B	211	LLP	C4-C4'-NZ	-2.22	113.80	124.04
1	D	211	LLP	OP3-P-OP2	2.19	116.02	107.80
1	B	211	LLP	O3-C3-C2	2.17	122.07	117.58
1	D	211	LLP	OP4-C5'-C5	2.16	113.40	109.36
1	C	211	LLP	C6-N1-C2	2.12	123.04	119.20
1	B	211	LLP	OP3-P-OP4	-2.07	101.28	106.67
1	D	211	LLP	C4-C4'-NZ	-2.03	114.69	124.04
1	A	211	LLP	C3-C4-C5	-2.02	116.65	118.28
1	B	211	LLP	OP3-P-OP2	2.01	115.32	107.80

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	211	LLP	O-C-CA-CB
1	B	211	LLP	O-C-CA-CB
1	C	211	LLP	O-C-CA-CB
1	D	211	LLP	O-C-CA-CB
1	B	211	LLP	C4-C4'-NZ-CE
1	A	211	LLP	C4-C4'-NZ-CE
1	C	211	LLP	C4-C4'-NZ-CE
1	D	211	LLP	CG-CD-CE-NZ
1	D	211	LLP	C4-C4'-NZ-CE
1	A	211	LLP	CD-CE-NZ-C4'
1	B	211	LLP	CD-CE-NZ-C4'
1	C	211	LLP	CG-CD-CE-NZ
1	D	211	LLP	CA-CB-CG-CD
1	B	211	LLP	C3-C4-C4'-NZ
1	A	211	LLP	C3-C4-C4'-NZ
1	C	211	LLP	C3-C4-C4'-NZ

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	391/398 (98%)	-0.01	25 (6%)	25 22	25, 44, 119, 148	0
1	B	391/398 (98%)	-0.32	13 (3%)	49 45	23, 35, 90, 131	0
1	C	395/398 (99%)	-0.19	21 (5%)	32 28	21, 35, 101, 144	0
1	D	391/398 (98%)	-0.26	14 (3%)	46 42	22, 38, 93, 128	0
All	All	1568/1592 (98%)	-0.19	73 (4%)	36 32	21, 37, 99, 148	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	SER	6.9
1	D	62	ILE	6.1
1	A	61	ARG	6.1
1	B	60	SER	5.5
1	D	59	TYR	5.5
1	A	58	PHE	5.0
1	A	59	TYR	5.0
1	D	58	PHE	5.0
1	C	365	ALA	4.8
1	A	62	ILE	4.7
1	D	367	TYR	4.6
1	C	51	ALA	4.6
1	A	55	ALA	4.5
1	A	57	HIS	4.5
1	A	51	ALA	4.4
1	A	365	ALA	4.3
1	D	61	ARG	4.3
1	B	63	SER	4.3
1	B	58	PHE	4.3
1	C	62	ILE	4.1
1	B	61	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	51	ALA	3.6
1	D	60	SER	3.5
1	C	60	SER	3.4
1	C	59	TYR	3.4
1	C	321	GLY	3.3
1	C	58	PHE	3.3
1	B	55	ALA	3.3
1	D	57	HIS	3.2
1	B	358	SER	3.2
1	D	63	SER	3.1
1	C	52	GLY	3.1
1	C	63	SER	3.1
1	C	367	TYR	3.1
1	A	368	GLY	3.0
1	B	59	TYR	2.9
1	C	55	ALA	2.9
1	A	52	GLY	2.9
1	A	54	GLN	2.8
1	B	48	ALA	2.8
1	A	367	TYR	2.8
1	C	50	PHE	2.8
1	B	57	HIS	2.8
1	A	359	TYR	2.7
1	B	44	GLU	2.6
1	D	56	GLY	2.6
1	C	322	ILE	2.6
1	C	366	HIS	2.6
1	C	3	GLY	2.5
1	C	47	ALA	2.5
1	B	52	GLY	2.5
1	C	45	TYR	2.5
1	A	361	PRO	2.5
1	B	359	TYR	2.4
1	B	62	ILE	2.4
1	A	355	THR	2.4
1	A	369	ILE	2.4
1	A	398	ALA	2.4
1	A	366	HIS	2.3
1	D	55	ALA	2.3
1	A	362	GLU	2.3
1	C	44	GLU	2.2
1	C	371	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	61	ARG	2.1
1	A	288	LEU	2.1
1	D	361	PRO	2.1
1	A	356	HIS	2.1
1	D	365	ALA	2.1
1	A	360	THR	2.1
1	C	41	PRO	2.1
1	D	360	THR	2.0
1	A	48	ALA	2.0
1	A	358	SER	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	LLP	B	211	24/25	0.97	0.06	25,30,32,33	0
1	LLP	C	211	24/25	0.97	0.06	24,26,27,28	0
1	LLP	A	211	24/25	0.98	0.05	30,32,33,36	0
1	LLP	D	211	24/25	0.98	0.05	26,29,31,32	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.