



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:51 PM UTC

PDB ID : 3P1T / pdb_00003p1t
Title : Crystal structure of a putative aminotransferase (BPSL1724) from Burkholderia pseudomallei K96243 at 2.60 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2010-09-30
Resolution : 2.60 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

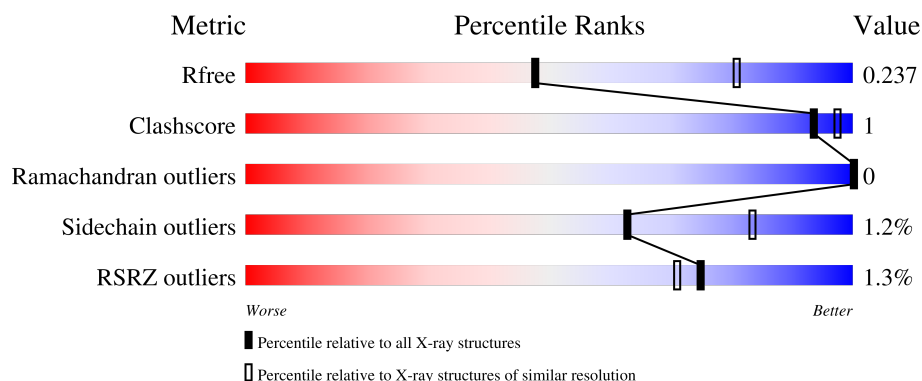
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	337	% 91% 5% .
1	B	337	% 90% 5% .
1	C	337	% 92% 5% .
1	D	337	% 88% 6% 6%

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	TLA	A	343	-	X	-	-
3	TLA	B	342	-	X	-	-
3	TLA	C	344	-	X	-	-
3	TLA	C	345	-	X	-	-
3	TLA	D	341	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10750 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 Putative histidinol-phosphate aminotransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	Se	0	3	0
			2512	1573	459	471	5	4			
1	B	319	Total	C	N	O	S	Se	0	4	0
			2515	1572	464	470	5	4			
1	C	321	Total	C	N	O	S	Se	0	4	0
			2530	1582	472	467	5	4			
1	D	317	Total	C	N	O	S	Se	0	2	0
			2492	1557	464	462	5	4			

There are 4 discrepancies between the modelled and reference sequences:

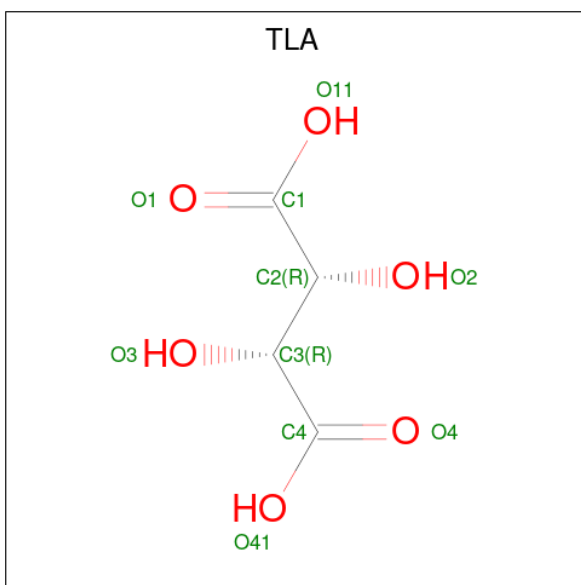
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q63U92
B	0	GLY	-	expression tag	UNP Q63U92
C	0	GLY	-	expression tag	UNP Q63U92
D	0	GLY	-	expression tag	UNP Q63U92

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



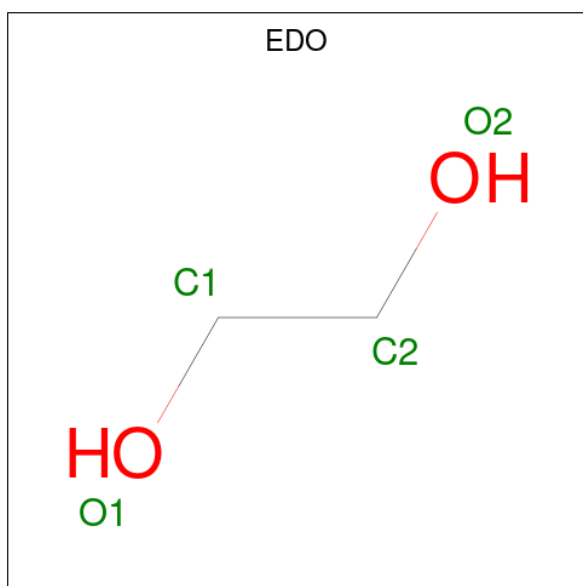
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is L(+)-TARTARIC ACID (CCD ID: TLA) (formula: C₄H₆O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 9 4 5	0	0
3	B	1	Total C O 10 4 6	0	0
3	C	1	Total C O 10 4 6	0	0
3	C	1	Total C O 10 4 6	0	0
3	D	1	Total C O 10 4 6	0	0

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 8 4 4	0	1

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 8 4 4	0	1
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 8 4 4	0	1

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	112	Total O 112 112	0	0
5	B	159	Total O 159 159	0	0
5	C	151	Total O 151 151	0	0
5	D	146	Total O 146 146	0	0

3 Residue-property plots [i](#)

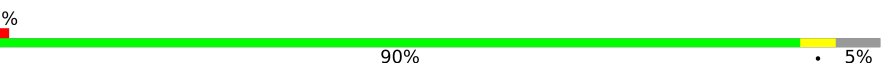
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Putative histidinol-phosphate aminotransferase

Chain A: 



- Molecule 1: Putative histidinol-phosphate aminotransferase

Chain B: 



- Molecule 1: Putative histidinol-phosphate aminotransferase

Chain C: 



- Molecule 1: Putative histidinol-phosphate aminotransferase

Chain D: 



HIS
SER

4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	167.55Å 167.55Å 290.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.65 – 2.60 29.65 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (29.65-2.60) 97.8 (29.65-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.61Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.8.0, BUSTER 2.8.0	Depositor
R, R_{free}	0.200 , 0.234 0.203 , 0.237	Depositor DCC
R_{free} test set	3131 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.378	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10750	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.49 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1359e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TLA, SO4, EDO

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	0.83	0/2561	1.23	3/3454 (0.1%)
1	B	0.85	0/2567	1.22	4/3460 (0.1%)
1	C	0.83	0/2582	1.21	4/3478 (0.1%)
1	D	0.86	1/2538 (0.0%)	1.22	4/3420 (0.1%)
All	All	0.84	1/10248 (0.0%)	1.22	15/13812 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	71	MSE	SE-CE	-5.61	1.78	1.95

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	321	ASP	CA-CB-CG	6.89	119.49	112.60
1	A	323	ASP	CA-CB-CG	6.41	119.01	112.60
1	B	18	VAL	CA-C-N	-5.94	115.11	122.84
1	B	18	VAL	C-N-CA	-5.94	115.11	122.84
1	A	135	ASP	CA-CB-CG	5.33	117.93	112.60
1	B	323	ASP	CA-CB-CG	5.19	117.79	112.60
1	D	302	ASP	CA-CB-CG	5.12	117.72	112.60
1	C	323	ASP	CA-CB-CG	5.10	117.70	112.60
1	D	180	ALA	CA-C-N	5.08	127.34	120.38
1	D	180	ALA	C-N-CA	5.08	127.34	120.38
1	C	153	ALA	CA-C-N	5.06	125.60	119.98
1	C	153	ALA	C-N-CA	5.06	125.60	119.98
1	D	77	ASP	CA-CB-CG	5.05	117.65	112.60
1	A	302	ASP	CA-CB-CG	5.01	117.61	112.60
1	B	77	ASP	CA-CB-CG	5.00	117.60	112.60

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	2512	0	2464	8	0
1	B	2515	0	2476	9	0
1	C	2530	0	2504	3	0
1	D	2492	0	2461	8	0
2	A	5	0	0	0	0
2	B	10	0	0	0	0
2	D	5	0	0	0	0
3	A	9	0	5	0	0
3	B	10	0	5	2	0
3	C	20	0	10	1	0
3	D	10	0	5	1	0
4	A	12	0	18	3	0
4	B	20	0	30	3	0
4	C	20	0	30	0	0
4	D	12	0	18	2	0
5	A	112	0	0	2	0
5	B	159	0	0	1	0
5	C	151	0	0	1	0
5	D	146	0	0	1	0
All	All	10750	0	10026	28	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 (28) 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:69:ASN:HD21	4:A:349:EDO:H12	1.46	0.81
1:B:43:ARG:NE	5:B:804:HOH:O	2.33	0.62
1:B:257:ARG:HH21	3:B:342:TLA:H3	1.65	0.62
1:D:274:SER:O	3:D:341:TLA:H3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:THR:HG21	4:B:353:EDO:H12	1.89	0.54
1:B:66:PRO:HD3	4:B:351:EDO:H22	1.90	0.53
1:D:81:ASP:OD1	1:D:105:ARG:HD3	2.10	0.52
1:B:81:ASP:OD1	1:B:105:ARG:HD3	2.10	0.52
1:A:84:SER:HB2	1:A:111:LEU:HD11	1.92	0.51
1:C:81:ASP:OD1	1:C:105:ARG:HD3	2.10	0.51
5:C:582:HOH:O	1:D:74:ARG:HD3	2.13	0.49
1:A:81:ASP:OD1	1:A:105:ARG:HD3	2.12	0.48
1:A:257:ARG:HE	4:A:346:EDO:H12	1.78	0.48
1:A:112:ARG:HD3	5:A:466:HOH:O	2.12	0.48
1:D:181:ARG:HD2	5:D:642:HOH:O	2.14	0.47
1:D:84:SER:HB2	1:D:111:LEU:HD11	1.96	0.47
1:A:215:LEU:HG	1:A:219:MSE:HE2	1.97	0.47
1:B:274:SER:O	3:B:342:TLA:H2	2.15	0.46
1:B:145:ASN:HD22	4:B:353:EDO:H22	1.82	0.45
1:C:87:PHE:HB3	1:C:90:MSE:SE	2.67	0.45
1:A:87:PHE:HB3	1:A:90:MSE:SE	2.67	0.45
1:B:26:PRO:HB2	1:B:318:ARG:HG3	1.99	0.43
1:D:99:PHE:CD1	4:D:354:EDO:H12	2.53	0.43
1:D:26:PRO:HB2	1:D:318:ARG:HG3	2.00	0.43
1:A:69:ASN:ND2	4:A:349:EDO:H12	2.26	0.42
5:A:438:HOH:O	1:B:74:ARG:HD3	2.20	0.41
1:D:145:ASN:HD22	4:D:354:EDO:H11	1.85	0.41
1:C:273:SER:HA	3:C:345:TLA:O3	2.21	0.41

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	323/337 (96%)	312 (97%)	11 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	321/337 (95%)	312 (97%)	9 (3%)	0	100	100
1	C	323/337 (96%)	313 (97%)	10 (3%)	0	100	100
1	D	317/337 (94%)	309 (98%)	8 (2%)	0	100	100
All	All	1284/1348 (95%)	1246 (97%)	38 (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	252/259 (97%)	248 (98%)	4 (2%)	55	79
1	B	255/259 (98%)	254 (100%)	1 (0%)	84	93
1	C	255/259 (98%)	251 (98%)	4 (2%)	55	79
1	D	253/259 (98%)	249 (98%)	4 (2%)	55	79
All	All	1015/1036 (98%)	1002 (99%)	13 (1%)	63	82

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	GLU
1	A	214[A]	GLU
1	A	214[B]	GLU
1	A	325	LEU
1	B	325	LEU
1	C	61	GLU
1	C	214	GLU
1	C	320	GLU
1	C	325	LEU
1	D	32	ARG
1	D	198	SER
1	D	320	GLU
1	D	325	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	227	ASN
1	A	311	HIS
1	B	127	ASN
1	B	227	ASN
1	B	283	ASN
1	B	299	GLN
1	C	69	ASN
1	C	223	GLN
1	C	283	ASN
1	D	223	GLN
1	D	227	ASN
1	D	283	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](#)

25 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	TLA	D	341	-	9,9,9	7.46	7 (77%)	12,12,12	3.74	8 (66%)
4	EDO	C	352	-	3,3,3	0.47	0	2,2,2	0.44	0
4	EDO	D	354	-	3,3,3	0.55	0	2,2,2	0.05	0
2	SO4	B	340	-	4,4,4	0.38	0	6,6,6	0.22	0
2	SO4	A	338	-	4,4,4	0.42	0	6,6,6	0.18	0
4	EDO	C	357[B]	-	3,3,3	0.55	0	2,2,2	0.24	0
3	TLA	A	343	-	6,8,9	6.76	5 (83%)	7,10,12	4.77	6 (85%)
4	EDO	D	358[A]	-	3,3,3	0.46	0	2,2,2	0.27	0
3	TLA	B	342	-	9,9,9	6.43	7 (77%)	12,12,12	3.55	6 (50%)
4	EDO	A	349	-	3,3,3	0.51	0	2,2,2	0.15	0
4	EDO	B	356[A]	-	3,3,3	0.50	0	2,2,2	0.08	0
3	TLA	C	344	-	9,9,9	7.08	7 (77%)	12,12,12	4.49	8 (66%)
2	SO4	B	339	-	4,4,4	0.35	0	6,6,6	0.16	0
4	EDO	C	357[A]	-	3,3,3	0.51	0	2,2,2	0.25	0
4	EDO	B	351	-	3,3,3	0.47	0	2,2,2	0.41	0
4	EDO	A	346	-	3,3,3	0.64	0	2,2,2	0.32	0
4	EDO	B	353	-	3,3,3	0.59	0	2,2,2	0.18	0
3	TLA	C	345	-	9,9,9	7.88	7 (77%)	12,12,12	3.69	7 (58%)
4	EDO	C	355	-	3,3,3	0.51	0	2,2,2	0.18	0
4	EDO	A	350	-	3,3,3	0.54	0	2,2,2	0.12	0
4	EDO	D	358[B]	-	3,3,3	0.43	0	2,2,2	0.41	0
4	EDO	B	356[B]	-	3,3,3	0.45	0	2,2,2	0.26	0
4	EDO	C	348	-	3,3,3	0.57	0	2,2,2	0.24	0
4	EDO	B	347	-	3,3,3	0.74	0	2,2,2	0.23	0
2	SO4	D	337	-	4,4,4	0.44	0	6,6,6	0.28	0

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	TLA	D	341	-	-	4/12/12/12	-
4	EDO	C	352	-	-	1/1/1/1	-
4	EDO	D	354	-	-	1/1/1/1	-
4	EDO	C	357[B]	-	-	0/1/1/1	-
3	TLA	A	343	-	-	3/9/10/12	-
4	EDO	D	358[A]	-	-	0/1/1/1	-
3	TLA	B	342	-	-	7/12/12/12	-
4	EDO	A	349	-	-	0/1/1/1	-
4	EDO	B	356[A]	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TLA	C	344	-	-	11/12/12/12	-
4	EDO	C	357[A]	-	-	0/1/1/1	-
4	EDO	B	351	-	-	0/1/1/1	-
4	EDO	A	346	-	-	0/1/1/1	-
4	EDO	B	353	-	-	0/1/1/1	-
3	TLA	C	345	-	-	7/12/12/12	-
4	EDO	C	355	-	-	0/1/1/1	-
4	EDO	A	350	-	-	0/1/1/1	-
4	EDO	D	358[B]	-	-	0/1/1/1	-
4	EDO	B	356[B]	-	-	0/1/1/1	-
4	EDO	C	348	-	-	0/1/1/1	-
4	EDO	B	347	-	-	0/1/1/1	-

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	345	TLA	C3-C4	14.12	1.72	1.52
3	D	341	TLA	C3-C4	12.26	1.70	1.52
3	C	345	TLA	C2-C1	10.55	1.67	1.52
3	C	344	TLA	O3-C3	10.53	1.62	1.42
3	C	345	TLA	O2-C2	9.89	1.61	1.42
3	B	342	TLA	O3-C3	9.80	1.61	1.42
3	D	341	TLA	O41-C4	9.04	1.59	1.30
3	C	344	TLA	C3-C4	8.99	1.65	1.52
3	A	343	TLA	O2-C2	8.93	1.59	1.43
3	D	341	TLA	O2-C2	8.91	1.59	1.42
3	D	341	TLA	O3-C3	8.90	1.59	1.42
3	C	344	TLA	O2-C2	8.85	1.59	1.42
3	B	342	TLA	C3-C4	8.84	1.65	1.52
3	C	344	TLA	O41-C4	8.70	1.58	1.30
3	A	343	TLA	O3-C3	8.27	1.58	1.42
3	C	345	TLA	O3-C3	7.88	1.57	1.42
3	A	343	TLA	C3-C4	7.87	1.63	1.52
3	B	342	TLA	O2-C2	7.73	1.57	1.42
3	B	342	TLA	O41-C4	7.72	1.55	1.30
3	C	344	TLA	C2-C1	7.62	1.63	1.52
3	D	341	TLA	C2-C1	7.37	1.63	1.52
3	A	343	TLA	O41-C4	6.38	1.51	1.30
3	C	345	TLA	C3-C2	6.24	1.72	1.53
3	C	345	TLA	O41-C4	5.81	1.49	1.30
3	B	342	TLA	C3-C2	5.47	1.70	1.53
3	D	341	TLA	O4-C4	5.46	1.38	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	341	TLA	C3-C2	5.12	1.69	1.53
3	B	342	TLA	C2-C1	5.09	1.59	1.52
3	C	344	TLA	C3-C2	4.80	1.68	1.53
3	A	343	TLA	O4-C4	4.76	1.36	1.22
3	C	344	TLA	O4-C4	4.73	1.36	1.22
3	B	342	TLA	O4-C4	4.55	1.35	1.22
3	C	345	TLA	O4-C4	3.46	1.32	1.22

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	344	TLA	O4-C4-C3	-8.65	98.59	121.62
3	D	341	TLA	O4-C4-C3	-8.10	100.04	121.62
3	B	342	TLA	O3-C3-C2	8.05	126.55	110.17
3	C	345	TLA	O2-C2-C3	7.59	125.62	110.17
3	A	343	TLA	O2-C2-C1	7.25	126.97	110.48
3	C	344	TLA	O2-C2-C1	7.14	125.94	110.69
3	D	341	TLA	O41-C4-C3	6.59	131.63	113.31
3	A	343	TLA	C3-C2-C1	-6.27	103.74	112.12
3	C	344	TLA	O3-C3-C2	6.17	122.72	110.17
3	C	344	TLA	O41-C4-C3	6.01	130.01	113.31
3	A	343	TLA	O4-C4-C3	-5.42	107.17	121.62
3	C	345	TLA	C3-C2-C1	-5.37	97.91	109.82
3	B	342	TLA	O4-C4-C3	-5.34	107.40	121.62
3	C	345	TLA	O41-C4-C3	5.04	127.30	113.31
3	B	342	TLA	O2-C2-C3	4.82	119.97	110.17
3	C	344	TLA	O3-C3-C4	4.45	120.20	110.69
3	D	341	TLA	C2-C3-C4	4.40	119.58	109.82
3	C	345	TLA	C2-C3-C4	4.17	119.06	109.82
3	A	343	TLA	O3-C3-C4	4.12	119.50	110.69
3	A	343	TLA	O41-C4-C3	3.65	123.44	113.31
3	B	342	TLA	O41-C4-C3	3.51	123.07	113.31
3	C	345	TLA	O3-C3-C2	3.24	116.77	110.17
3	C	344	TLA	O41-C4-O4	3.22	131.39	124.08
3	D	341	TLA	O3-C3-C2	3.17	116.62	110.17
3	C	345	TLA	O4-C4-C3	-3.05	113.48	121.62
3	D	341	TLA	C3-C2-C1	2.85	116.14	109.82
3	B	342	TLA	O2-C2-C1	2.84	116.77	110.69
3	D	341	TLA	O11-C1-C2	2.71	120.84	113.31
3	D	341	TLA	O1-C1-C2	-2.51	114.93	121.62
3	C	345	TLA	O41-C4-O4	-2.44	118.54	124.08
3	C	344	TLA	C3-C2-C1	-2.40	104.50	109.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	342	TLA	O41-C4-O4	2.40	129.52	124.08
3	A	343	TLA	O41-C4-O4	2.33	129.36	124.08
3	C	344	TLA	C2-C3-C4	-2.15	105.06	109.82
3	D	341	TLA	O2-C2-C1	2.06	115.09	110.69

There are no chirality outliers.

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	343	TLA	C1-C2-C3-O3
3	A	343	TLA	C1-C2-C3-C4
3	A	343	TLA	O2-C2-C3-C4
3	C	344	TLA	C2-C3-C4-O4
3	C	344	TLA	C2-C3-C4-O41
3	C	345	TLA	O1-C1-C2-O2
3	C	345	TLA	O11-C1-C2-O2
3	D	341	TLA	C1-C2-C3-C4
3	C	344	TLA	C1-C2-C3-C4
3	B	342	TLA	O11-C1-C2-C3
3	B	342	TLA	C2-C3-C4-O4
3	B	342	TLA	C2-C3-C4-O41
3	B	342	TLA	O2-C2-C3-O3
3	C	345	TLA	C1-C2-C3-C4
3	B	342	TLA	O3-C3-C4-O4
3	B	342	TLA	O3-C3-C4-O41
3	B	342	TLA	O1-C1-C2-C3
3	C	345	TLA	O1-C1-C2-C3
3	C	345	TLA	O2-C2-C3-C4
3	C	345	TLA	O11-C1-C2-C3
3	D	341	TLA	C1-C2-C3-O3
3	D	341	TLA	O2-C2-C3-C4
3	C	344	TLA	O11-C1-C2-O2
3	C	344	TLA	O2-C2-C3-C4
3	C	344	TLA	O1-C1-C2-O2
3	D	341	TLA	O11-C1-C2-O2
3	C	344	TLA	O3-C3-C4-O41
3	C	344	TLA	O3-C3-C4-O4
3	C	345	TLA	O3-C3-C4-O41
4	C	352	EDO	O1-C1-C2-O2
3	C	344	TLA	O11-C1-C2-C3
4	D	354	EDO	O1-C1-C2-O2
3	C	344	TLA	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	C	344	TLA	C1-C2-C3-O3

There are no ring outliers.

8 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	341	TLA	1	0
4	D	354	EDO	2	0
3	B	342	TLA	2	0
4	A	349	EDO	2	0
4	B	351	EDO	1	0
4	A	346	EDO	1	0
4	B	353	EDO	2	0
3	C	345	TLA	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	318/337 (94%)	0.01	5 (1%) 70 66	29, 46, 72, 114	3 (0%)
1	B	315/337 (93%)	-0.21	3 (0%) 79 76	24, 39, 55, 94	4 (1%)
1	C	317/337 (94%)	-0.21	5 (1%) 70 66	24, 39, 62, 104	4 (1%)
1	D	313/337 (92%)	-0.27	3 (0%) 79 76	22, 37, 58, 75	2 (0%)
All	All	1263/1348 (93%)	-0.17	16 (1%) 75 71	22, 40, 63, 114	13 (1%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	335	HIS	6.8
1	A	14	ALA	3.8
1	D	333	SER	3.5
1	A	334	ASP	3.1
1	C	221[A]	ARG	3.0
1	C	14	ALA	3.0
1	C	334	ASP	2.9
1	B	244[A]	ARG	2.9
1	B	17	ALA	2.8
1	A	13	ALA	2.8
1	A	318	ARG	2.4
1	C	43[A]	ARG	2.3
1	D	17	ALA	2.2
1	C	184	ALA	2.1
1	A	15	ALA	2.1
1	D	241	ASP	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	EDO	C	357[A]	4/4	0.56	0.29	34,36,36,40	4
4	EDO	C	357[B]	4/4	0.56	0.29	31,31,33,34	4
4	EDO	D	358[A]	4/4	0.58	0.29	47,48,49,50	4
4	EDO	D	358[B]	4/4	0.58	0.29	52,53,55,59	4
4	EDO	B	356[A]	4/4	0.69	0.18	46,46,48,51	4
4	EDO	B	356[B]	4/4	0.69	0.18	73,75,76,78	4
4	EDO	D	354	4/4	0.70	0.36	72,72,73,79	0
4	EDO	A	350	4/4	0.72	0.21	58,60,60,64	0
4	EDO	A	346	4/4	0.73	0.19	44,48,48,51	0
4	EDO	A	349	4/4	0.73	0.20	64,66,67,69	0
4	EDO	B	351	4/4	0.76	0.17	59,60,63,64	0
3	TLA	C	345	10/10	0.77	0.21	57,69,72,73	0
4	EDO	B	347	4/4	0.79	0.17	41,42,43,47	0
2	SO4	A	338	5/5	0.79	0.25	100,105,105,107	0
4	EDO	C	352	4/4	0.79	0.19	59,61,62,64	0
4	EDO	B	353	4/4	0.79	0.35	71,71,75,77	0
3	TLA	C	344	10/10	0.80	0.16	57,65,73,73	0
3	TLA	D	341	10/10	0.80	0.23	75,81,92,93	0
4	EDO	C	355	4/4	0.81	0.25	74,75,75,80	0
3	TLA	B	342	10/10	0.83	0.20	77,81,87,89	0
2	SO4	B	339	5/5	0.84	0.15	93,97,99,99	0
2	SO4	B	340	5/5	0.86	0.23	47,51,53,53	5
2	SO4	D	337	5/5	0.87	0.14	91,96,97,97	0
3	TLA	A	343	9/10	0.88	0.17	68,73,77,80	0
4	EDO	C	348	4/4	0.89	0.16	44,45,45,46	0

6.5 Other polymers ⓘ

There are no such residues in this entry.