



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:37 PM UTC

PDB ID : 3ORW / pdb_00003orw
Title : Crystal structure of thermophilic phosphotriesterase from *Geobacillus kaustophilus* HTA426
Authors : Zheng, B.S.; Yu, S.S.; Zhang, Y.; Lou, Z.Y.; Feng, Y.
Deposited on : 2010-09-07
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

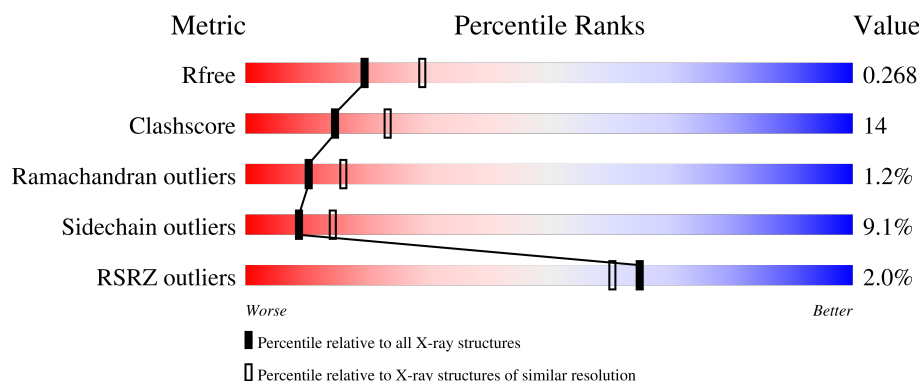
1 Overall quality at a glance

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	326	% 68% 25% 5% ..
1	B	326	3% 56% 34% 7% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5101 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 Phosphotriesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2548	1620	436	476	16			
1	B	324	Total	C	N	O	S	0	0	0
			2548	1620	436	476	16			

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Co	0	0
			2	2		
2	B	2	Total	Co	0	0
			2	2		

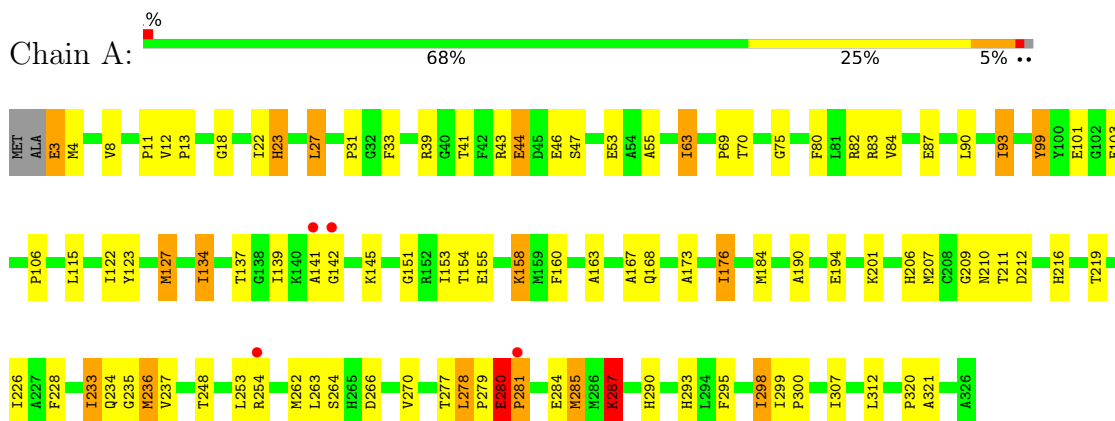
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	O	0	0
			1	1		

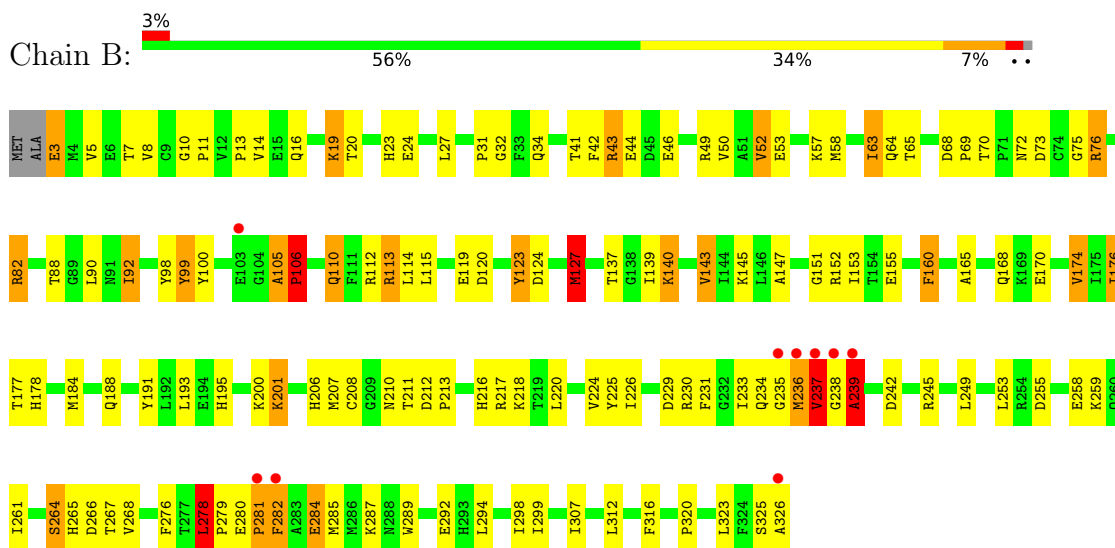
3 Residue-property plots

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: Phosphotriesterase



• Molecule 1: Phosphotriesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.32Å 88.88Å 89.38Å 90.00° 98.58° 90.00°	Depositor
Resolution (Å)	47.15 – 2.40 47.15 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.2 (47.15-2.40) 96.2 (47.15-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.196 , 0.268 0.199 , 0.268	Depositor DCC
R_{free} test set	1502 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 22.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5101	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: KCX, CO

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	1.67	18/2594 (0.7%)	1.50	23/3510 (0.7%)
1	B	1.75	36/2594 (1.4%)	1.56	30/3510 (0.9%)
All	All	1.71	54/5188 (1.0%)	1.53	53/7020 (0.8%)

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	2

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	299	ILE	CA-CB	10.38	1.59	1.54
1	B	106	PRO	CA-C	9.93	1.61	1.52
1	B	307	ILE	CA-CB	8.90	1.64	1.54
1	A	63	ILE	CA-CB	8.55	1.64	1.54
1	B	165	ALA	CA-CB	-7.49	1.41	1.53
1	B	299	ILE	CA-CB	7.48	1.58	1.54
1	B	268	VAL	CA-CB	7.33	1.62	1.53
1	A	163	ALA	CA-CB	-7.19	1.42	1.53
1	A	233	ILE	CA-CB	6.97	1.61	1.53
1	B	52	VAL	N-CA	6.94	1.54	1.46
1	B	63	ILE	CA-CB	6.52	1.61	1.54
1	B	105	ALA	CA-CB	6.44	1.60	1.53
1	B	239	ALA	CA-C	6.37	1.60	1.53
1	B	123	TYR	CA-C	6.37	1.61	1.52
1	B	174	VAL	C-O	6.35	1.32	1.24
1	B	174	VAL	CA-C	6.27	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	174	VAL	CA-CB	6.24	1.61	1.54
1	A	84	VAL	CA-CB	6.15	1.62	1.54
1	A	69	PRO	CA-C	6.13	1.61	1.52
1	A	321	ALA	CA-CB	6.12	1.63	1.53
1	B	239	ALA	N-CA	6.06	1.54	1.45
1	B	42	PHE	N-CA	-6.01	1.38	1.46
1	B	237	VAL	N-CA	5.97	1.53	1.46
1	B	143	VAL	CA-CB	5.89	1.65	1.55
1	B	224	VAL	CA-CB	5.88	1.62	1.54
1	B	73	ASP	N-CA	5.85	1.53	1.46
1	A	70	THR	C-N	5.82	1.41	1.33
1	A	206	HIS	C-O	5.80	1.31	1.24
1	B	70	THR	C-O	5.76	1.32	1.24
1	A	75	GLY	C-O	-5.72	1.16	1.24
1	A	190	ALA	CA-CB	5.62	1.62	1.53
1	B	88	THR	CA-CB	5.61	1.62	1.54
1	B	52	VAL	CA-CB	5.61	1.60	1.54
1	A	127	MET	N-CA	5.55	1.53	1.46
1	B	233	ILE	CA-CB	5.46	1.60	1.53
1	B	65	THR	C-O	5.46	1.29	1.23
1	B	19	LYS	CD-CE	5.42	1.68	1.52
1	B	237	VAL	CA-CB	5.39	1.61	1.54
1	A	122	ILE	CA-CB	5.39	1.60	1.54
1	A	219	THR	CA-CB	5.35	1.62	1.53
1	A	18	GLY	N-CA	5.34	1.53	1.45
1	B	160	PHE	CA-C	5.34	1.59	1.52
1	B	19	LYS	N-CA	5.33	1.52	1.46
1	B	92	ILE	CA-CB	5.25	1.60	1.54
1	B	120	ASP	CA-C	5.25	1.59	1.52
1	B	225	TYR	N-CA	-5.24	1.39	1.45
1	B	299	ILE	CA-C	5.23	1.57	1.52
1	A	141	ALA	CA-CB	5.21	1.61	1.53
1	A	264	SER	C-O	5.20	1.29	1.24
1	B	105	ALA	CA-C	5.20	1.58	1.53
1	B	170	GLU	C-O	-5.14	1.17	1.24
1	A	134	ILE	C-O	5.10	1.30	1.24
1	B	119	GLU	N-CA	-5.06	1.40	1.46
1	B	153	ILE	CA-CB	5.03	1.59	1.53

All (53) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	PRO	N-CA-C	11.37	124.57	110.70
1	B	280	GLU	CA-C-N	-11.11	106.74	118.85
1	B	280	GLU	C-N-CA	-11.11	106.74	118.85
1	A	33	PHE	N-CA-C	-8.75	101.67	111.82
1	B	52	VAL	N-CA-CB	8.49	120.48	110.55
1	A	160	PHE	N-CA-C	-7.93	101.94	111.69
1	B	177	THR	N-CA-C	7.75	121.10	109.95
1	B	52	VAL	N-CA-C	7.71	117.82	110.42
1	B	106	PRO	N-CA-C	7.42	119.75	110.70
1	A	219	THR	N-CA-C	-7.34	102.66	111.69
1	B	278	LEU	CA-C-N	7.04	128.65	119.84
1	B	278	LEU	C-N-CA	7.04	128.65	119.84
1	A	93	ILE	N-CA-C	-7.00	98.38	108.53
1	B	76	ARG	NE-CZ-NH1	-6.87	114.63	121.50
1	B	140	LYS	N-CA-C	-6.82	99.19	110.17
1	B	41	THR	N-CA-C	6.63	119.84	110.23
1	B	208	CYS	N-CA-C	6.59	120.42	112.38
1	B	231	PHE	N-CA-C	6.52	119.82	110.24
1	A	278	LEU	CA-C-N	6.49	127.51	119.98
1	A	278	LEU	C-N-CA	6.49	127.51	119.98
1	A	158	LYS	N-CA-C	6.21	118.04	111.28
1	B	50	VAL	CB-CA-C	-6.21	103.66	112.22
1	A	44	GLU	N-CA-C	6.10	117.62	110.97
1	B	261	ILE	N-CA-C	6.03	116.55	108.11
1	A	298	ILE	N-CA-CB	5.85	117.26	110.65
1	A	293	HIS	N-CA-C	-5.82	104.89	112.23
1	B	7	THR	N-CA-C	-5.79	100.61	109.41
1	A	153	ILE	N-CA-C	-5.79	98.09	106.88
1	A	115	LEU	N-CA-C	-5.75	106.10	113.23
1	B	280	GLU	N-CA-C	5.75	122.51	109.81
1	B	298	ILE	N-CA-C	5.68	116.33	110.82
1	B	127	MET	CG-SD-CE	-5.63	88.51	100.90
1	A	280	GLU	N-CA-C	5.62	122.22	109.81
1	A	209	GLY	N-CA-C	-5.59	108.00	115.32
1	A	212	ASP	CA-C-N	5.55	125.22	119.56
1	A	212	ASP	C-N-CA	5.55	125.22	119.56
1	B	70	THR	CA-C-N	5.39	125.32	119.76
1	B	70	THR	C-N-CA	5.39	125.32	119.76
1	B	105	ALA	CA-C-N	-5.39	114.82	120.38
1	B	105	ALA	C-N-CA	-5.39	114.82	120.38
1	A	176	ILE	CA-C-N	-5.36	113.46	123.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	176	ILE	C-N-CA	-5.36	113.46	123.05
1	B	225	TYR	N-CA-C	-5.21	102.81	110.52
1	A	290	HIS	N-CA-C	-5.20	101.16	109.07
1	B	281	PRO	N-CA-C	-5.10	106.56	113.40
1	B	10	GLY	CA-C-N	5.10	124.86	119.76
1	B	10	GLY	C-N-CA	5.10	124.86	119.76
1	B	292	GLU	CA-C-N	5.09	127.41	120.54
1	B	292	GLU	C-N-CA	5.09	127.41	120.54
1	A	101	GLU	CA-C-N	5.03	125.68	119.99
1	A	101	GLU	C-N-CA	5.03	125.68	119.99
1	B	127	MET	N-CA-C	5.02	117.40	111.33
1	A	43	ARG	CG-CD-NE	-5.01	100.98	112.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	234	GLN	Peptide
1	B	239	ALA	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	2548	0	2492	61	0
1	B	2548	0	2492	80	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	B	1	0	0	0	0
All	All	5101	0	4984	140	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 14.

All (140) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:MET:HE2	1:B:282:PHE:CZ	1.70	1.26
1:A:134:ILE:O	1:A:137:THR:HG22	1.08	1.26
1:A:134:ILE:O	1:A:137:THR:CG2	1.87	1.23
1:A:235:GLY:HA3	1:A:285:MET:HE1	1.38	1.05
1:A:151:GLY:HA2	1:A:184:MET:HE3	1.40	1.02
1:A:234:GLN:O	1:A:285:MET:HE3	1.63	0.99
1:A:262:MET:HE3	1:A:320:PRO:HG3	1.46	0.98
1:B:236:MET:HE2	1:B:282:PHE:CE2	2.00	0.96
1:A:127:MET:CE	1:A:167:ALA:HA	2.08	0.84
1:A:123:TYR:CE1	1:A:127:MET:HE3	2.12	0.83
1:B:235:GLY:HA3	1:B:285:MET:SD	2.19	0.81
1:A:137:THR:HG23	1:A:139:ILE:H	1.47	0.79
1:B:325:SER:O	1:B:326:ALA:HB2	1.83	0.78
1:B:43:ARG:HB2	1:B:43:ARG:HH11	1.49	0.78
1:B:112:ARG:HA	1:B:115:LEU:HD12	1.65	0.76
1:B:210:ASN:O	1:B:216:HIS:HE1	1.68	0.76
1:B:325:SER:O	1:B:326:ALA:CB	2.34	0.76
1:A:235:GLY:HA3	1:A:285:MET:CE	2.15	0.76
1:A:262:MET:CE	1:A:320:PRO:HG3	2.14	0.76
1:A:4:MET:HE3	1:A:11:PRO:HB2	1.68	0.76
1:A:127:MET:HE2	1:A:167:ALA:HA	1.66	0.75
1:A:168:GLN:HG3	1:A:173:ALA:O	1.86	0.75
1:A:262:MET:HE3	1:A:320:PRO:CG	2.17	0.73
1:B:110:GLN:HE21	1:B:110:GLN:HA	1.54	0.73
1:B:32:GLY:H	1:B:34:GLN:NE2	1.86	0.73
1:B:236:MET:HE2	1:B:282:PHE:HZ	1.48	0.73
1:B:282:PHE:HA	1:B:285:MET:HE3	1.71	0.72
1:B:32:GLY:H	1:B:34:GLN:HE21	1.36	0.71
1:B:13:PRO:HG2	1:B:16:GLN:NE2	2.04	0.71
1:A:123:TYR:HE1	1:A:127:MET:HE3	1.53	0.71
1:B:123:TYR:CZ	1:B:127:MET:HE3	2.25	0.71
1:B:265:HIS:HB2	1:B:267:THR:HG23	1.75	0.69
1:B:113:ARG:CZ	1:B:155:GLU:OE2	2.41	0.69
1:A:207:MET:HG3	1:A:226:ILE:HB	1.76	0.68
1:B:123:TYR:CE1	1:B:127:MET:CE	2.77	0.68
1:A:123:TYR:CE1	1:A:127:MET:CE	2.78	0.66
1:B:210:ASN:O	1:B:216:HIS:CE1	2.49	0.66
1:A:55:ALA:HB1	1:A:90:LEU:HD22	1.78	0.65
1:A:253:LEU:HD13	1:A:307:ILE:HD12	1.79	0.64
1:B:43:ARG:HB2	1:B:43:ARG:NH1	2.12	0.64
1:B:137:THR:HB	1:B:139:ILE:HD12	1.80	0.64
1:A:123:TYR:HE1	1:A:127:MET:CE	2.11	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ASN:O	1:A:216:HIS:HE1	1.82	0.62
1:B:123:TYR:CE1	1:B:127:MET:HE1	2.34	0.62
1:B:236:MET:CE	1:B:282:PHE:CE2	2.82	0.61
1:A:137:THR:HG23	1:A:139:ILE:N	2.14	0.61
1:B:235:GLY:CA	1:B:285:MET:SD	2.89	0.60
1:B:13:PRO:HG2	1:B:16:GLN:HE21	1.63	0.60
1:B:123:TYR:CE1	1:B:127:MET:HE3	2.37	0.59
1:B:237:VAL:O	1:B:237:VAL:HG12	2.03	0.59
1:B:284:GLU:O	1:B:287:LYS:HG2	2.02	0.59
1:A:155:GLU:OE1	1:A:158:LYS:NZ	2.36	0.59
1:B:220:LEU:HD21	1:B:226:ILE:HD13	1.86	0.58
1:A:127:MET:HE2	1:A:167:ALA:CB	2.34	0.57
1:B:191:TYR:O	1:B:195:HIS:HD2	1.88	0.57
1:B:68:ASP:OD1	1:B:68:ASP:C	2.47	0.56
1:B:242:ASP:OD1	1:B:245:ARG:NH1	2.39	0.56
1:A:151:GLY:CA	1:A:184:MET:HE3	2.25	0.56
1:A:127:MET:HE1	1:A:167:ALA:HA	1.87	0.55
1:B:281:PRO:HB2	1:B:285:MET:HE2	1.88	0.55
1:B:113:ARG:NH2	1:B:155:GLU:CD	2.66	0.54
1:B:19:LYS:HB3	1:B:63:ILE:HD13	1.90	0.54
1:A:47:SER:HB2	1:A:270:VAL:HG11	1.89	0.54
1:A:99:TYR:HD2	1:A:103:GLU:O	1.92	0.53
1:A:210:ASN:O	1:A:216:HIS:CE1	2.61	0.53
1:B:113:ARG:NH2	1:B:155:GLU:OE2	2.42	0.53
1:B:184:MET:O	1:B:188:GLN:HG3	2.08	0.53
1:B:236:MET:HG3	1:B:239:ALA:HB3	1.90	0.53
1:A:284:GLU:O	1:A:287:LYS:HB2	2.09	0.53
1:B:100:TYR:HB3	1:B:147:ALA:HB1	1.90	0.53
1:B:113:ARG:HG2	1:B:113:ARG:NH1	2.24	0.53
1:B:58:MET:O	1:B:63:ILE:HB	2.09	0.52
1:B:98:TYR:HB2	1:B:160:PHE:CE2	2.44	0.52
1:B:82:ARG:HH11	1:B:82:ARG:HG2	1.75	0.51
1:A:127:MET:HE2	1:A:167:ALA:CA	2.36	0.51
1:A:280:GLU:OE2	1:A:280:GLU:HA	2.10	0.51
1:A:280:GLU:N	1:A:281:PRO:HD2	2.26	0.51
1:B:106:PRO:O	1:B:110:GLN:HG2	2.11	0.50
1:B:229:ASP:O	1:B:266:ASP:HB2	2.12	0.50
1:B:220:LEU:HD21	1:B:226:ILE:CD1	2.40	0.50
1:A:134:ILE:HB	1:A:137:THR:HG21	1.93	0.49
1:A:137:THR:OG1	1:A:139:ILE:HD12	2.12	0.49
1:A:39:ARG:HE	1:B:124:ASP:CG	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:GLU:HA	1:A:46:GLU:OE1	2.12	0.49
1:B:127:MET:HE2	1:B:127:MET:HB2	1.38	0.49
1:A:63:ILE:HD11	1:A:295:PHE:CZ	2.48	0.49
1:B:43:ARG:HH12	1:B:46:GLU:HB3	1.78	0.49
1:A:312:LEU:HD23	1:A:312:LEU:HA	1.65	0.48
1:B:174:VAL:HG13	1:B:323:LEU:HD11	1.94	0.48
1:B:75:GLY:O	1:B:76:ARG:C	2.56	0.48
1:A:55:ALA:CB	1:A:90:LEU:HD22	2.41	0.48
1:A:3:GLU:HG2	1:A:4:MET:HG2	1.97	0.47
1:B:143:VAL:HG11	1:B:176:ILE:HD13	1.97	0.47
1:B:276:PHE:HE2	1:B:278:LEU:HD21	1.80	0.47
1:B:43:ARG:NH1	1:B:46:GLU:HB3	2.30	0.47
1:B:23:HIS:CD2	1:B:69:PRO:HG3	2.50	0.46
1:A:228:PHE:HB2	1:A:263:LEU:HD23	1.98	0.46
1:A:4:MET:HE3	1:A:11:PRO:CB	2.42	0.46
1:A:236:MET:O	1:A:237:VAL:HG12	2.16	0.46
1:A:237:VAL:O	1:A:237:VAL:HG13	2.15	0.45
1:A:87:GLU:O	1:A:87:GLU:HG3	2.15	0.45
1:B:212:ASP:OD1	1:B:212:ASP:C	2.60	0.45
1:A:8:VAL:O	1:A:82:ARG:HG3	2.17	0.45
1:B:207:MET:O	1:B:216:HIS:NE2	2.49	0.45
1:B:200:LYS:C	1:B:201:LYS:HG2	2.40	0.45
1:B:206:HIS:CG	1:B:230:ARG:HE	2.35	0.44
1:A:44:GLU:HG2	1:A:83:ARG:HH21	1.83	0.44
1:B:8:VAL:HG13	1:B:92:ILE:O	2.18	0.44
1:A:93:ILE:HG23	1:A:142:GLY:HA3	2.00	0.43
1:B:253:LEU:C	1:B:255:ASP:H	2.25	0.43
1:B:113:ARG:NE	1:B:155:GLU:OE2	2.51	0.43
1:A:280:GLU:OE2	1:A:280:GLU:CA	2.66	0.43
1:B:207:MET:HG3	1:B:226:ILE:HB	2.01	0.42
1:B:312:LEU:O	1:B:316:PHE:HD1	2.02	0.42
1:A:262:MET:HE3	1:A:320:PRO:CD	2.48	0.42
1:A:154:THR:O	1:A:158:LYS:HG3	2.19	0.42
1:A:27:LEU:HD21	1:A:80:PHE:CE2	2.55	0.42
1:B:211:THR:O	1:B:213:PRO:HD3	2.20	0.42
1:B:24:GLU:OE2	1:B:264:SER:OG	2.31	0.42
1:A:216:HIS:CE1	1:A:248:THR:HG21	2.55	0.41
1:A:278:LEU:HA	1:A:279:PRO:HD2	1.81	0.41
1:B:31:PRO:HB3	1:B:105:ALA:HB2	2.01	0.41
1:B:32:GLY:N	1:B:34:GLN:HE21	2.10	0.41
1:A:298:ILE:HD13	1:A:298:ILE:HA	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:ILE:HD13	1:A:262:MET:HE1	2.03	0.41
1:B:3:GLU:O	1:B:14:VAL:HG23	2.21	0.41
1:A:23:HIS:CD2	1:A:266:ASP:OD1	2.73	0.41
1:B:23:HIS:CD2	1:B:69:PRO:CG	3.04	0.41
1:B:53:GLU:OE1	1:B:57:LYS:HE3	2.20	0.41
1:B:259:LYS:HZ3	1:B:259:LYS:HG3	1.67	0.41
1:B:34:GLN:CD	1:B:34:GLN:H	2.28	0.41
1:B:99:TYR:HD1	1:B:99:TYR:HA	1.80	0.41
1:A:279:PRO:C	1:A:281:PRO:HD2	2.46	0.41
1:B:151:GLY:C	1:B:184:MET:HE3	2.46	0.41
1:B:20:THR:HB	1:B:320:PRO:HG2	2.02	0.41
1:B:5:VAL:O	1:B:11:PRO:HA	2.21	0.40
1:B:64:GLN:O	1:B:90:LEU:HD12	2.21	0.40
1:B:147:ALA:HA	1:B:178:HIS:HB3	2.04	0.40
1:A:12:VAL:HG13	1:A:13:PRO:HD2	2.03	0.40
1:B:191:TYR:O	1:B:195:HIS:CD2	2.72	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	321/326 (98%)	302 (94%)	15 (5%)	4 (1%)	10	16
1	B	321/326 (98%)	301 (94%)	16 (5%)	4 (1%)	10	16
All	All	642/652 (98%)	603 (94%)	31 (5%)	8 (1%)	10	16

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	281	PRO
1	B	237	VAL

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Mol	Chain	Res	Type
1	B	279	PRO
1	A	287	LYS
1	B	238	GLY
1	B	289	TRP
1	A	23	HIS
1	A	280	GLU

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	263/264 (100%)	245 (93%)	18 (7%)	14	25
1	B	263/264 (100%)	233 (89%)	30 (11%)	5	8
All	All	526/528 (100%)	478 (91%)	48 (9%)	9	14

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	27	LEU
1	A	31	PRO
1	A	41	THR
1	A	53	GLU
1	A	99	TYR
1	A	176	ILE
1	A	194	GLU
1	A	201	LYS
1	A	211	THR
1	A	233	ILE
1	A	236	MET
1	A	254	ARG
1	A	277	THR
1	A	280	GLU
1	A	285	MET
1	A	287	LYS

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Mol	Chain	Res	Type
1	A	300	PRO
1	B	3	GLU
1	B	27	LEU
1	B	43	ARG
1	B	44	GLU
1	B	49	ARG
1	B	52	VAL
1	B	72	ASN
1	B	82	ARG
1	B	99	TYR
1	B	106	PRO
1	B	110	GLN
1	B	113	ARG
1	B	114	LEU
1	B	127	MET
1	B	140	LYS
1	B	152	ARG
1	B	168	GLN
1	B	176	ILE
1	B	193	LEU
1	B	201	LYS
1	B	217	ARG
1	B	218	LYS
1	B	236	MET
1	B	249	LEU
1	B	258	GLU
1	B	264	SER
1	B	278	LEU
1	B	282	PHE
1	B	284	GLU
1	B	294	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	91	ASN
1	A	195	HIS
1	A	216	HIS
1	B	16	GLN
1	B	34	GLN
1	B	110	GLN

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Mol	Chain	Res	Type
1	B	168	GLN
1	B	195	HIS
1	B	216	HIS
1	B	290	HIS
1	B	297	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	KCX	A	145	2,1	10,11,12	2.36	2 (20%)	6,12,14	0.83	0
1	KCX	B	145	2,1	10,11,12	2.52	3 (30%)	6,12,14	1.38	1 (16%)

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	KCX	A	145	2,1	-	1/9/10/12	-
1	KCX	B	145	2,1	-	1/9/10/12	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	145	KCX	OQ1-CX	5.54	1.31	1.21
1	A	145	KCX	OQ1-CX	5.36	1.31	1.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	145	KCX	O-C	4.95	1.38	1.20
1	A	145	KCX	O-C	4.81	1.38	1.20
1	B	145	KCX	CX-NZ	2.12	1.39	1.35

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	145	KCX	OQ1-CX-NZ	-2.22	121.56	124.92

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	145	KCX	CA-CB-CG-CD
1	A	145	KCX	CA-CB-CG-CD

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](#)

Of 4 ligands modelled in this entry, 4 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 ⓘ

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	323/326 (99%)	-0.14	4 (1%) 76 73	25, 38, 55, 68	0
1	B	323/326 (99%)	0.03	9 (2%) 55 51	26, 40, 59, 82	0
All	All	646/652 (99%)	-0.05	13 (2%) 65 60	25, 39, 57, 82	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	239	ALA	7.9
1	B	237	VAL	5.7
1	B	235	GLY	5.0
1	B	282	PHE	4.9
1	B	326	ALA	4.7
1	B	238	GLY	3.6
1	B	236	MET	3.4
1	A	281	PRO	2.9
1	B	281	PRO	2.6
1	A	142	GLY	2.3
1	B	103	GLU	2.3
1	A	141	ALA	2.2
1	A	254	ARG	2.1

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	KCX	B	145	12/13	0.95	0.08	28,29,40,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	KCX	A	145	12/13	0.96	0.09	29,32,35,37	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CO	A	328	1/1	0.97	0.04	43,43,43,43	0
2	CO	B	328	1/1	0.98	0.04	54,54,54,54	0
2	CO	B	327	1/1	0.99	0.05	48,48,48,48	0
2	CO	A	327	1/1	0.99	0.02	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.