



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:46 AM GMT

PDB ID : 2O6I
Title : Structure of an Enterococcus Faecalis HD Domain Phosphohydrolase
Authors : Vorontsov, I.I.; Minasov, G.; Shuvalova, L.; Brunzelle, J.S.; Moy, S.; Col-lart, F.R.; Joachimiak, A.; Anderson, W.F.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2006-12-07
Resolution : 2.55 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

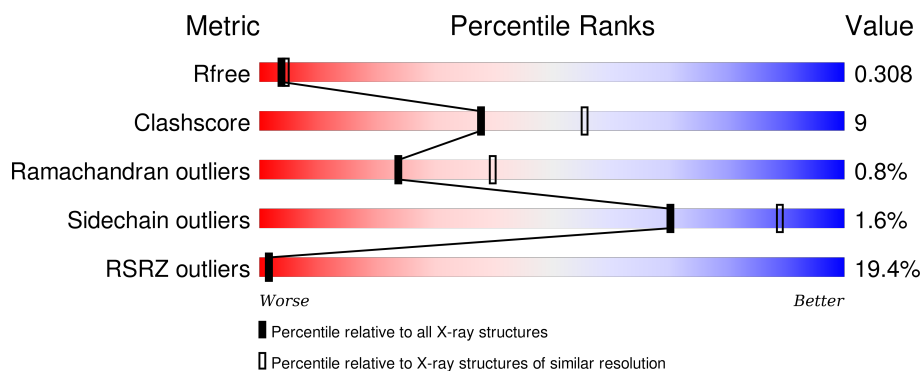
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	91344	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	480	
1	B	480	

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	CL	A	503	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7556 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 HD domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	436	Total	C	N	O	S	0	6	0
			3632	2325	618	676	13			
1	B	430	Total	C	N	O	S	0	12	0
			3610	2308	614	675	13			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	INITIATING METHIONINE	UNP Q836G9
A	-22	HIS	-	EXPRESSION TAG	UNP Q836G9
A	-21	HIS	-	EXPRESSION TAG	UNP Q836G9
A	-20	HIS	-	EXPRESSION TAG	UNP Q836G9
A	-19	HIS	-	EXPRESSION TAG	UNP Q836G9
A	-18	HIS	-	EXPRESSION TAG	UNP Q836G9
A	-17	HIS	-	EXPRESSION TAG	UNP Q836G9
A	-16	SER	-	CLONING ARTIFACT	UNP Q836G9
A	-15	SER	-	CLONING ARTIFACT	UNP Q836G9
A	-14	GLY	-	CLONING ARTIFACT	UNP Q836G9
A	-13	VAL	-	CLONING ARTIFACT	UNP Q836G9
A	-12	ASP	-	CLONING ARTIFACT	UNP Q836G9
A	-11	LEU	-	CLONING ARTIFACT	UNP Q836G9
A	-10	GLY	-	CLONING ARTIFACT	UNP Q836G9
A	-9	THR	-	CLONING ARTIFACT	UNP Q836G9
A	-8	GLU	-	CLONING ARTIFACT	UNP Q836G9
A	-7	ASN	-	CLONING ARTIFACT	UNP Q836G9
A	-6	LEU	-	CLONING ARTIFACT	UNP Q836G9
A	-5	TYR	-	CLONING ARTIFACT	UNP Q836G9
A	-4	PHE	-	CLONING ARTIFACT	UNP Q836G9
A	-3	GLN	-	CLONING ARTIFACT	UNP Q836G9
A	-2	SER	-	CLONING ARTIFACT	UNP Q836G9
A	-1	ASN	-	CLONING ARTIFACT	UNP Q836G9
A	0	ALA	-	CLONING ARTIFACT	UNP Q836G9
B	-23	MET	-	CLONING ARTIFACT	UNP Q836G9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	HIS	-	EXPRESSION TAG	UNP Q836G9
B	-21	HIS	-	EXPRESSION TAG	UNP Q836G9
B	-20	HIS	-	EXPRESSION TAG	UNP Q836G9
B	-19	HIS	-	EXPRESSION TAG	UNP Q836G9
B	-18	HIS	-	EXPRESSION TAG	UNP Q836G9
B	-17	HIS	-	EXPRESSION TAG	UNP Q836G9
B	-16	SER	-	CLONING ARTIFACT	UNP Q836G9
B	-15	SER	-	CLONING ARTIFACT	UNP Q836G9
B	-14	GLY	-	CLONING ARTIFACT	UNP Q836G9
B	-13	VAL	-	CLONING ARTIFACT	UNP Q836G9
B	-12	ASP	-	CLONING ARTIFACT	UNP Q836G9
B	-11	LEU	-	CLONING ARTIFACT	UNP Q836G9
B	-10	GLY	-	CLONING ARTIFACT	UNP Q836G9
B	-9	THR	-	CLONING ARTIFACT	UNP Q836G9
B	-8	GLU	-	CLONING ARTIFACT	UNP Q836G9
B	-7	ASN	-	CLONING ARTIFACT	UNP Q836G9
B	-6	LEU	-	CLONING ARTIFACT	UNP Q836G9
B	-5	TYR	-	CLONING ARTIFACT	UNP Q836G9
B	-4	PHE	-	CLONING ARTIFACT	UNP Q836G9
B	-3	GLN	-	CLONING ARTIFACT	UNP Q836G9
B	-2	SER	-	CLONING ARTIFACT	UNP Q836G9
B	-1	ASN	-	CLONING ARTIFACT	UNP Q836G9
B	0	ALA	-	CLONING ARTIFACT	UNP Q836G9

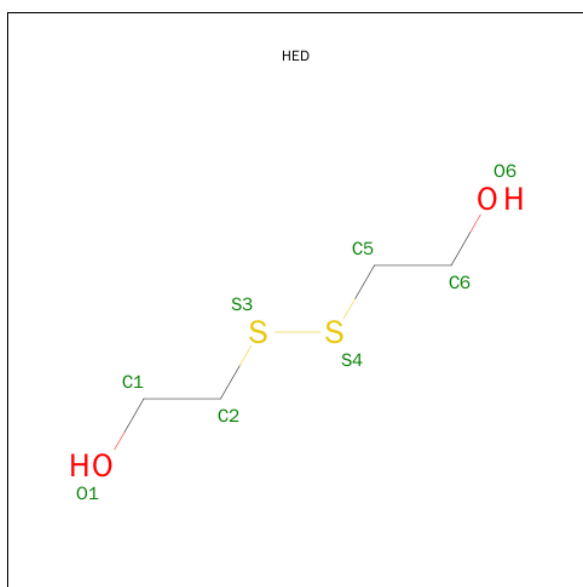
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	1
3	A	1	Total Cl 1 1	0	0

- Molecule 4 is 2-HYDROXYETHYL DISULFIDE (three-letter code: HED) (formula: C₄H₁₀O₂S₂).



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

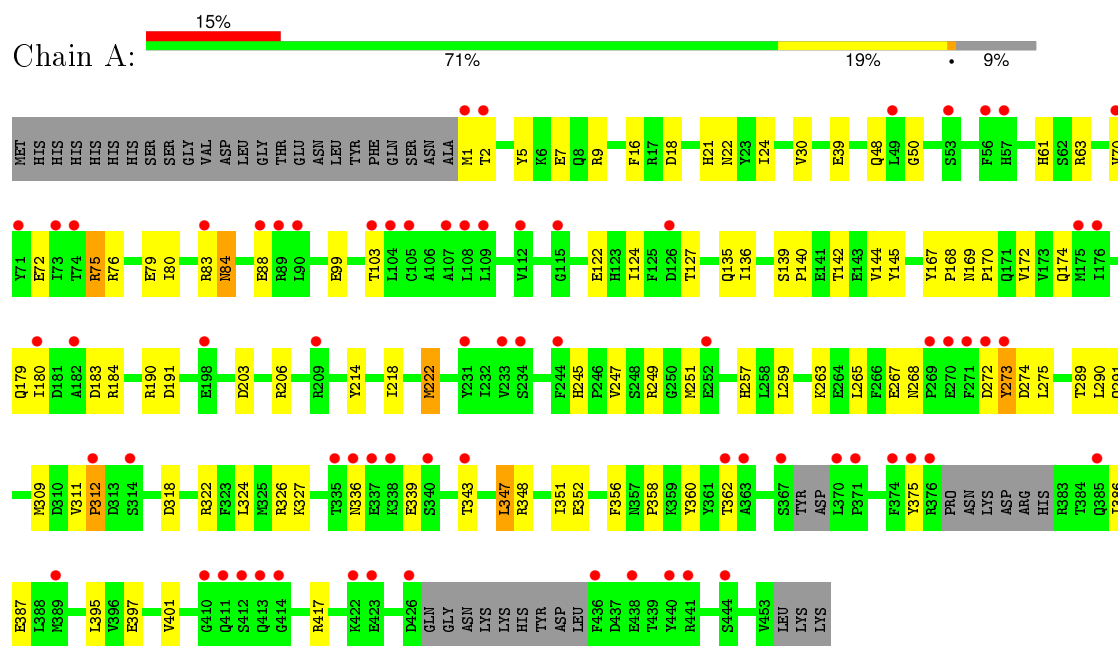
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	169	Total	O	0	2
			170	170		
5	B	128	Total	O	0	4
			132	132		

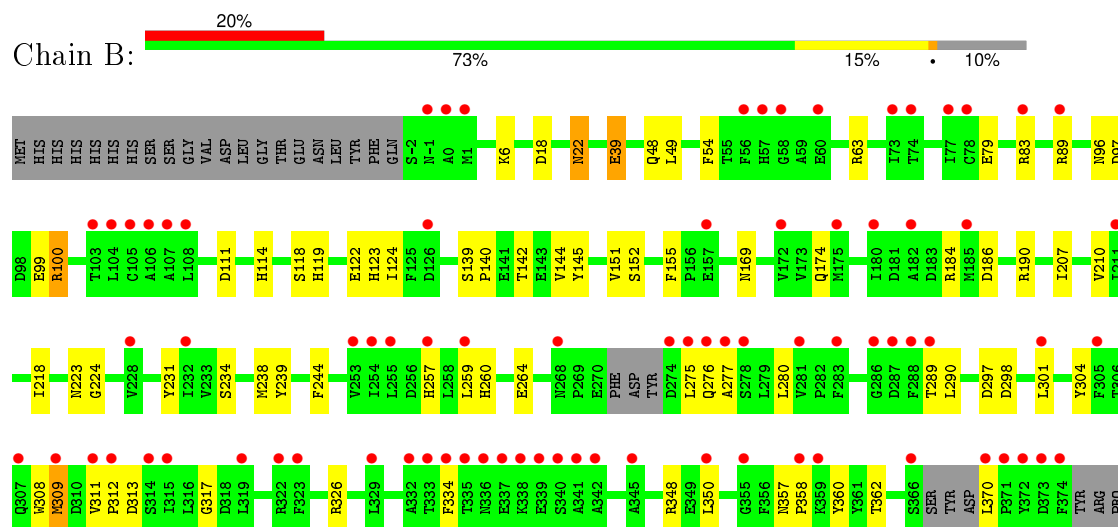
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: HD domain protein



• Molecule 1: HD domain protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.91Å 109.91Å 182.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.55 29.97 – 2.55	Depositor EDS
% Data completeness (in resolution range)	96.3 (30.00-2.55) 96.3 (29.97-2.55)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.248 , 0.317 0.241 , 0.308	Depositor DCC
R_{free} test set	2049 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	57.5	Xtriage
Anisotropy	0.001	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 65.5	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 40601 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7556	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: HED, ZN, CL

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	0.29	0/3718	0.54	0/5032
1	B	0.28	0/3694	0.53	1/5000 (0.0%)
All	All	0.28	0/7412	0.54	1/10032 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	49	LEU	CA-CB-CG	5.54	128.03	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3632	0	3542	76	0
1	B	3610	0	3512	56	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	8	0	10	1	0
5	A	170	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	132	0	0	0	0
All	All	7556	0	7064	132	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 9.

All (132) 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:75:ARG:HG3	1:A:75:ARG:HH11	1.25	1.01
1:A:75:ARG:CG	1:A:75:ARG:HH11	1.85	0.88
1:B:18:ASP:O	1:B:22:ASN:HA	1.75	0.86
1:B:257:HIS:HD2	1:B:360:TYR:HA	1.46	0.81
1:A:265:LEU:HD13	1:A:272:ASP:OD2	1.81	0.80
1:B:100:ARG:CG	1:B:100:ARG:HH11	2.01	0.73
1:A:16:PHE:HB2	1:A:24:ILE:HB	1.72	0.71
1:A:75:ARG:HG3	1:A:75:ARG:NH1	1.94	0.69
1:A:124:ILE:HD13	1:A:259:LEU:HB2	1.72	0.69
1:B:111:ASP:HB2	1:B:114:HIS:ND1	2.08	0.68
1:A:249:ARG:NH2	1:A:417:ARG:HG3	2.10	0.66
1:A:80:ILE:O	1:A:84:ASN:ND2	2.29	0.66
1:A:1:MET:N	1:A:2:THR:HA	2.12	0.65
1:B:39:GLU:CD	1:B:39:GLU:H	2.01	0.64
1:A:375:TYR:HB3	1:A:386:ILE:HD11	1.81	0.63
1:B:297:ASP:OD2	1:B:298:ASP:N	2.33	0.61
1:A:140:PRO:HA	1:A:145:TYR:CD2	2.35	0.61
1:B:118:SER:O	1:B:122:GLU:HB2	2.02	0.60
1:A:124:ILE:CD1	1:A:259:LEU:HB2	2.31	0.60
1:A:79:GLU:O	1:A:83:ARG:HG2	2.02	0.59
1:A:257:HIS:HD2	1:A:360:TYR:HA	1.66	0.59
1:B:100:ARG:HH11	1:B:100:ARG:HG2	1.68	0.59
1:A:351:ILE:HG22	1:A:356:PHE:HB2	1.86	0.58
1:A:245:HIS:CD2	1:A:247:VAL:HG22	2.39	0.58
1:A:88:GLU:OE2	1:A:88:GLU:N	2.36	0.57
1:B:18:ASP:O	1:B:22:ASN:CA	2.49	0.57
1:B:99:GLU:OE2	1:B:169:ASN:ND2	2.38	0.57
1:B:275:LEU:O	1:B:277:ALA:N	2.37	0.57
1:A:386:ILE:CG2	1:A:397:GLU:HB2	2.35	0.57
1:A:21:HIS:O	1:A:22:ASN:HB2	2.05	0.56
1:B:111:ASP:HB2	1:B:114:HIS:CE1	2.40	0.56
1:B:152:SER:HB3	1:B:155:PHE:HB2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:ASP:O	1:A:322:ARG:HG3	2.06	0.56
1:A:142:THR:HG22	1:A:144:VAL:H	1.71	0.56
1:A:7:GLU:O	1:A:9[B]:ARG:HD2	2.06	0.55
1:A:18:ASP:OD1	1:A:190:ARG:NH2	2.39	0.55
1:B:54:PHE:HD1	1:B:326:ARG:HG2	1.71	0.55
1:A:103:THR:HB	1:A:172:VAL:HG13	1.88	0.55
1:A:214:TYR:CE1	1:A:395:LEU:HD11	2.42	0.55
1:A:358:PRO:O	1:A:362:THR:HB	2.07	0.55
1:B:184:ARG:HH22	1:B:370:LEU:HD13	1.70	0.55
1:A:72:GLU:O	1:A:76:ARG:HG2	2.07	0.54
1:B:239:TYR:HD2	1:B:244:PHE:CE2	2.25	0.54
1:B:210[B]:VAL:HG11	1:B:224:GLY:HA3	1.90	0.53
1:A:251:MET:HE3	1:A:326:ARG:HA	1.91	0.52
1:A:251:MET:HE1	1:A:326:ARG:C	2.30	0.52
1:B:124:ILE:CD1	1:B:259:LEU:HB2	2.39	0.52
1:B:124:ILE:HD13	1:B:259:LEU:HB2	1.91	0.51
1:A:122:GLU:HA	1:A:127[A]:THR:HG23	1.93	0.51
1:A:84:ASN:HD22	1:A:84:ASN:H	1.60	0.50
1:B:234:SER:O	1:B:238:MET:HB2	2.11	0.50
1:B:100:ARG:NH1	1:B:100:ARG:CG	2.68	0.50
1:A:84:ASN:HD22	1:A:84:ASN:N	2.08	0.50
1:B:6:LYS:HB3	1:B:151:VAL:HG13	1.94	0.50
1:A:5:TYR:CE2	1:A:30:VAL:HG22	2.47	0.49
1:B:142:THR:HG22	1:B:144:VAL:H	1.77	0.49
1:A:348:ARG:HE	1:A:362:THR:HG21	1.77	0.49
1:A:135:GLN:OE1	4:A:505:HED:H21	2.12	0.49
1:A:142:THR:HG22	1:A:144:VAL:N	2.28	0.48
1:A:48:GLN:HE21	1:A:63:ARG:HA	1.79	0.48
1:B:100:ARG:HG3	1:B:100:ARG:HH11	1.79	0.48
1:A:386:ILE:HG21	1:A:397:GLU:HB2	1.96	0.47
1:B:275:LEU:HD23	1:B:277:ALA:H	1.79	0.47
1:B:79:GLU:HG3	1:B:83[A]:ARG:HE	1.79	0.47
1:B:97:ASP:OD1	1:B:100:ARG:NH1	2.47	0.47
1:B:309:MET:HA	1:B:317:GLY:HA2	1.97	0.46
1:A:61:HIS:CD2	1:A:61:HIS:H	2.32	0.46
1:B:48:GLN:NE2	1:B:63:ARG:HG2	2.31	0.46
1:B:257:HIS:CD2	1:B:360:TYR:HA	2.36	0.46
1:A:289:THR:HG22	1:A:291:GLN:H	1.81	0.46
1:A:352:GLU:HG2	1:A:358:PRO:HD3	1.98	0.46
1:A:174:GLN:NE2	1:A:218:ILE:H	2.14	0.46
1:B:260:HIS:O	1:B:264:GLU:HG3	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:TYR:CG	1:A:274:ASP:N	2.84	0.45
1:B:142:THR:HG21	1:B:144:VAL:HG22	1.98	0.45
1:A:139:SER:HA	1:A:140:PRO:HD3	1.75	0.45
1:A:99:GLU:O	1:A:103:THR:HG22	2.17	0.45
1:A:289:THR:HG22	1:A:290:LEU:N	2.30	0.45
1:A:309:MET:CE	1:A:324:LEU:HD12	2.47	0.45
1:A:122:GLU:HG3	1:A:127[B]:THR:O	2.17	0.44
1:A:48:GLN:HG2	5:A:569:HOH:O	2.18	0.44
1:A:84:ASN:N	1:A:84:ASN:ND2	2.65	0.44
1:A:347:LEU:O	1:A:351:ILE:HG13	2.18	0.44
1:B:139:SER:HA	1:B:140:PRO:HD3	1.84	0.44
1:A:311:VAL:HA	1:A:312:PRO:HD2	1.85	0.44
1:A:336:ASN:HB3	1:A:339:GLU:HB2	1.98	0.44
1:B:39:GLU:OE2	1:B:142:THR:HG23	2.17	0.44
1:B:140:PRO:HA	1:B:145:TYR:CD2	2.53	0.44
1:B:334:PHE:HB3	1:B:452:LEU:HD12	1.99	0.44
1:B:142:THR:CG2	1:B:144:VAL:HG22	2.48	0.44
1:B:48:GLN:HE21	1:B:63:ARG:HA	1.83	0.44
1:B:18:ASP:O	1:B:22:ASN:N	2.51	0.43
1:B:357:ASN:HA	1:B:358:PRO:HD2	1.90	0.43
1:B:348:ARG:HG2	1:B:362:THR:HG21	2.00	0.43
1:A:75:ARG:CG	1:A:75:ARG:NH1	2.55	0.43
1:A:251:MET:CE	1:A:326:ARG:HA	2.47	0.43
1:B:437:ASP:HA	1:B:440:TYR:HD1	1.83	0.43
1:B:289:THR:HG22	1:B:290:LEU:N	2.32	0.43
1:A:263:LYS:NZ	1:A:267:GLU:OE2	2.42	0.43
1:A:124:ILE:HD13	1:A:259:LEU:CB	2.45	0.43
1:A:136:ILE:HG23	1:A:142:THR:HG21	2.01	0.43
1:B:174:GLN:NE2	1:B:218:ILE:H	2.16	0.43
1:B:207:ILE:HD11	1:B:231:TYR:HB2	2.01	0.43
1:A:169:ASN:HA	1:A:170:PRO:HD2	1.88	0.43
1:B:79:GLU:HG3	1:B:83[B]:ARG:HE	1.83	0.42
1:A:257:HIS:CD2	1:A:360:TYR:HA	2.50	0.42
1:A:343:THR:O	1:A:347:LEU:HG	2.18	0.42
1:B:119[B]:HIS:O	1:B:123[B]:HIS:ND1	2.49	0.42
1:A:61:HIS:HE1	1:A:191:ASP:OD1	2.01	0.42
1:A:309:MET:HE3	1:A:324:LEU:HD12	2.00	0.42
1:B:280:LEU:HD21	1:B:301:LEU:HD21	2.00	0.42
1:A:179:GLN:HA	1:A:184:ARG:NH1	2.34	0.42
1:A:351:ILE:CG2	1:A:356:PHE:HB2	2.48	0.42
1:B:350:LEU:HB3	1:B:439:THR:HG23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:GLU:OE1	1:B:100:ARG:NH2	2.53	0.42
1:B:304:TYR:O	1:B:308:TRP:CD1	2.73	0.42
1:B:311:VAL:HG12	1:B:313:ASP:H	1.85	0.41
1:A:203:ASP:OD2	1:A:206:ARG:HG3	2.21	0.41
1:A:70:VAL:HG22	1:A:183:ASP:HA	2.02	0.41
1:A:222:MET:HE1	1:A:401:VAL:HG11	2.03	0.41
1:A:48:GLN:HA	1:A:63:ARG:HG2	2.03	0.41
1:B:223[A]:ASN:H	1:B:223[A]:ASN:HD22	1.69	0.41
1:B:142:THR:HG22	1:B:144:VAL:N	2.35	0.41
1:B:275:LEU:C	1:B:277:ALA:H	2.24	0.41
1:A:336:ASN:CB	1:A:339:GLU:HB2	2.51	0.41
1:A:122:GLU:HG3	1:A:127[A]:THR:O	2.21	0.41
1:A:401:VAL:O	1:A:401:VAL:HG12	2.20	0.41
1:A:39:GLU:CD	1:A:142:THR:HG23	2.41	0.40
1:A:322:ARG:HG2	1:A:327:LYS:HE3	2.04	0.40
1:A:167:TYR:HA	1:A:168:PRO:HD3	1.91	0.40
1:A:249:ARG:NH2	1:A:417:ARG:CG	2.81	0.40
1:B:186:ASP:OD1	1:B:190:ARG:NH1	2.54	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	434/480 (90%)	405 (93%)	25 (6%)	4 (1%)	21	36
1	B	432/480 (90%)	406 (94%)	23 (5%)	3 (1%)	26	44
All	All	866/960 (90%)	811 (94%)	48 (6%)	7 (1%)	24	40

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	276	GLN
1	A	50	GLY
1	A	273	TYR
1	B	22	ASN
1	A	180	ILE
1	B	312	PRO
1	A	312	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	396/430 (92%)	389 (98%)	7 (2%)	66	87
1	B	395/430 (92%)	390 (99%)	5 (1%)	76	91
All	All	791/860 (92%)	779 (98%)	12 (2%)	70	90

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	ARG
1	A	84	ASN
1	A	222	MET
1	A	268	ASN
1	A	275	LEU
1	A	347	LEU
1	A	387	GLU
1	B	39	GLU
1	B	89	ARG
1	B	96	ASN
1	B	100	ARG
1	B	309	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	25	HIS
1	A	27	GLN
1	A	48	GLN
1	A	61	HIS
1	A	84	ASN
1	A	114	HIS
1	A	149	ASN
1	A	174	GLN
1	A	223	ASN
1	A	245	HIS
1	A	257	HIS
1	A	268	ASN
1	A	276	GLN
1	A	307	GLN
1	A	336	ASN
1	A	390	GLN
1	A	411	GLN
1	A	413	GLN
1	B	8	GLN
1	B	48	GLN
1	B	61	HIS
1	B	96	ASN
1	B	128	ASN
1	B	146	GLN
1	B	149	ASN
1	B	174	GLN
1	B	237	GLN
1	B	241	GLN
1	B	268	ASN
1	B	385	GLN
1	B	413	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
4	HED	A	505	-	7,7,7	0.64	0	6,6,6	1.22	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
4	HED	A	505	-	-	0/5/5/5	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	505	HED	1	0

5.7 Other polymers

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

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	436/480 (90%)	1.02	70 (16%) 3 2	57, 67, 78, 96	0
1	B	430/480 (89%)	1.31	98 (22%) 1 1	57, 69, 81, 101	0
All	All	866/960 (90%)	1.16	168 (19%) 1 1	57, 68, 79, 101	0

All (168) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	376	ARG	8.0
1	A	414	GLY	7.9
1	B	58	GLY	7.3
1	B	371	PRO	7.2
1	A	271	PHE	7.0
1	A	374	PHE	6.9
1	B	412	SER	6.7
1	B	312	PRO	6.6
1	A	337	GLU	6.3
1	B	335	THR	6.2
1	B	338	LYS	6.0
1	B	374	PHE	5.8
1	A	338	LYS	5.7
1	B	287	ASP	5.6
1	A	335	THR	5.6
1	B	423	GLU	5.6
1	B	370	LEU	5.5
1	B	289	THR	5.4
1	A	273	TYR	5.4
1	B	337	GLU	5.4
1	B	355	GLY	5.4
1	A	270	GLU	5.4
1	B	275	LEU	5.3
1	B	336	ASN	5.2

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Mol	Chain	Res	Type	RSRZ
1	B	373	ASP	5.0
1	B	413	GLN	4.7
1	B	386	ILE	4.7
1	B	410	GLY	4.6
1	A	411	GLN	4.5
1	A	336	ASN	4.5
1	A	126	ASP	4.4
1	B	439	THR	4.3
1	B	411	GLN	4.2
1	B	274	ASP	4.2
1	A	108	LEU	4.2
1	A	441	ARG	4.1
1	B	126	ASP	4.1
1	A	272	ASP	4.0
1	A	410	GLY	4.0
1	A	413	GLN	4.0
1	B	255	LEU	3.9
1	B	414	GLY	3.9
1	A	370	LEU	3.9
1	B	449	ASN	3.8
1	B	440	TYR	3.8
1	B	268	ASN	3.8
1	B	175	MET	3.7
1	B	366	SER	3.7
1	B	57	HIS	3.7
1	B	286	GLY	3.7
1	B	334	PHE	3.5
1	B	276	GLN	3.5
1	A	423	GLU	3.5
1	A	74	THR	3.5
1	A	107	ALA	3.5
1	B	283	PHE	3.4
1	B	446	TYR	3.4
1	A	176	ILE	3.4
1	A	104	LEU	3.4
1	A	371	PRO	3.4
1	A	57	HIS	3.4
1	B	372	TYR	3.4
1	A	2	THR	3.3
1	B	389	MET	3.2
1	B	309	MET	3.2
1	A	89	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	415	ASP	3.2
1	A	1	MET	3.2
1	A	367	SER	3.1
1	B	339	GLU	3.1
1	B	107	ALA	3.1
1	A	175	MET	3.1
1	B	60	GLU	3.1
1	A	70	VAL	3.1
1	A	182	ALA	3.0
1	A	180	ILE	3.0
1	B	182	ALA	3.0
1	B	277	ALA	3.0
1	B	323	PHE	3.0
1	A	103	THR	3.0
1	B	301	LEU	2.9
1	B	288	PHE	2.9
1	A	269	PRO	2.9
1	B	333	THR	2.9
1	B	1	MET	2.9
1	B	74	THR	2.9
1	B	305	PHE	2.9
1	B	311	VAL	2.8
1	B	104	LEU	2.8
1	A	73	ILE	2.8
1	A	314	SER	2.8
1	A	375	TYR	2.8
1	B	253	VAL	2.8
1	B	232	ILE	2.8
1	B	185	MET	2.8
1	A	88	GLU	2.7
1	A	440	TYR	2.7
1	B	329	LEU	2.7
1	A	56	PHE	2.7
1	B	89	ARG	2.7
1	A	362	THR	2.7
1	B	315	ILE	2.7
1	B	341	ALA	2.7
1	A	115	GLY	2.7
1	A	422	LYS	2.7
1	B	314	SER	2.6
1	B	-1	ASN	2.6
1	B	259	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	112	VAL	2.6
1	A	363	ALA	2.6
1	A	252[A]	GLU	2.6
1	A	244	PHE	2.6
1	B	108	LEU	2.6
1	A	53	SER	2.6
1	B	278	SER	2.6
1	A	412	SER	2.5
1	B	322	ARG	2.5
1	B	56	PHE	2.5
1	B	445	SER	2.5
1	A	385	GLN	2.5
1	B	228	VAL	2.5
1	B	358	PRO	2.5
1	A	343	THR	2.5
1	A	109	LEU	2.4
1	B	425	LEU	2.4
1	A	209	ARG	2.4
1	B	436	PHE	2.4
1	B	254	ILE	2.4
1	B	257	HIS	2.4
1	B	157	GLU	2.4
1	A	90	LEU	2.4
1	A	105	CYS	2.4
1	A	71	TYR	2.4
1	B	281	VAL	2.4
1	A	340	SER	2.4
1	B	105	CYS	2.3
1	B	437	ASP	2.3
1	A	83	ARG	2.3
1	B	307	GLN	2.3
1	B	340	SER	2.3
1	A	426	ASP	2.3
1	B	211	ILE	2.3
1	B	180	ILE	2.3
1	B	350	LEU	2.3
1	A	436	PHE	2.2
1	B	106	ALA	2.2
1	B	424	MET	2.2
1	A	198	GLU	2.2
1	B	103	THR	2.2
1	B	172	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	332	ALA	2.2
1	B	345	ALA	2.2
1	B	73	ILE	2.2
1	A	231	TYR	2.1
1	B	83[A]	ARG	2.1
1	B	0	ALA	2.1
1	B	359	LYS	2.1
1	A	49	LEU	2.1
1	B	77	ILE	2.1
1	A	234	SER	2.1
1	A	438	GLU	2.1
1	B	78	CYS	2.1
1	B	342	ALA	2.1
1	A	389	MET	2.0
1	A	312	PRO	2.0
1	A	444	SER	2.0
1	A	233	VAL	2.0
1	B	319	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	CL	A	503	1/1	0.81	0.41	10.73	84,84,84,84	1
3	CL	B	504[A]	1/1	0.96	0.17	-1.36	63,63,63,63	1
2	ZN	A	501	1/1	0.91	0.11	-2.53	68,68,68,68	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	B	502	1/1	0.98	0.09	-4.37	74,74,74,74	1
4	HED	A	505	8/8	0.91	0.18	-	63,64,66,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.