



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 02:50 AM BST

PDB ID : 3ZLF
Title : Structure of group A Streptococcal enolase K312A mutant
Authors : Cork, A.J.; Ericsson, D.J.; Law, R.H.P.; Casey, L.W.; Valkov, E.; Bertozzi, C.; Stamp, A.; Aquilina, J.A.; Whisstock, J.C.; Walker, M.J.; Kobe, B.
Deposited on : 2013-01-31
Resolution : 2.15 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

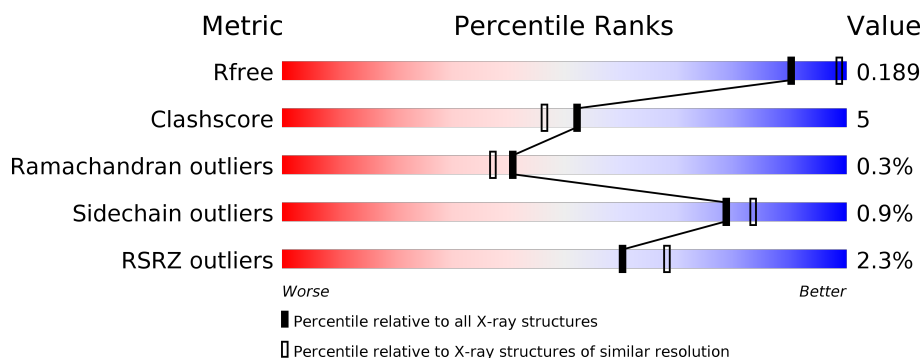
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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}	130704	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 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	455	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 11%, green 85%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 85% 11% • </div> </div>
1	B	455	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 1%, yellow 11%, green 84%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 84% 11% 5% </div> </div>
1	C	455	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 83%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 83% 12% • </div> </div>
1	D	455	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 4%, orange 1%, yellow 11%, green 84%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 4% 84% 11% • 5% </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	437	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 27040 atoms, of which 13121 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 ENOLASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	436	Total	C	H	N	O	S	0	0	0
			6636	2102	3298	562	663	11			
1	B	434	Total	C	H	N	O	S	0	0	0
			6592	2090	3272	558	661	11			
1	C	435	Total	C	H	N	O	S	0	0	0
			6603	2093	3279	558	662	11			
1	D	434	Total	C	H	N	O	S	0	0	0
			6592	2090	3272	558	661	11			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q5XD01
A	-18	GLY	-	expression tag	UNP Q5XD01
A	-17	SER	-	expression tag	UNP Q5XD01
A	-16	SER	-	expression tag	UNP Q5XD01
A	-15	HIS	-	expression tag	UNP Q5XD01
A	-14	HIS	-	expression tag	UNP Q5XD01
A	-13	HIS	-	expression tag	UNP Q5XD01
A	-12	HIS	-	expression tag	UNP Q5XD01
A	-11	HIS	-	expression tag	UNP Q5XD01
A	-10	HIS	-	expression tag	UNP Q5XD01
A	-9	SER	-	expression tag	UNP Q5XD01
A	-8	SER	-	expression tag	UNP Q5XD01
A	-7	GLY	-	expression tag	UNP Q5XD01
A	-6	LEU	-	expression tag	UNP Q5XD01
A	-5	VAL	-	expression tag	UNP Q5XD01
A	-4	PRO	-	expression tag	UNP Q5XD01
A	-3	ARG	-	expression tag	UNP Q5XD01
A	-2	GLY	-	expression tag	UNP Q5XD01
A	-1	SER	-	expression tag	UNP Q5XD01
A	0	HIS	-	expression tag	UNP Q5XD01
A	312	ALA	LYS	engineered mutation	UNP Q5XD01

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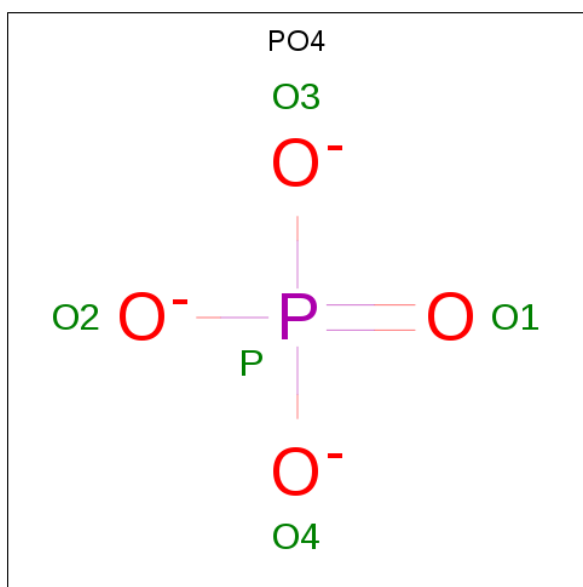
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	expression tag	UNP Q5XD01
B	-18	GLY	-	expression tag	UNP Q5XD01
B	-17	SER	-	expression tag	UNP Q5XD01
B	-16	SER	-	expression tag	UNP Q5XD01
B	-15	HIS	-	expression tag	UNP Q5XD01
B	-14	HIS	-	expression tag	UNP Q5XD01
B	-13	HIS	-	expression tag	UNP Q5XD01
B	-12	HIS	-	expression tag	UNP Q5XD01
B	-11	HIS	-	expression tag	UNP Q5XD01
B	-10	HIS	-	expression tag	UNP Q5XD01
B	-9	SER	-	expression tag	UNP Q5XD01
B	-8	SER	-	expression tag	UNP Q5XD01
B	-7	GLY	-	expression tag	UNP Q5XD01
B	-6	LEU	-	expression tag	UNP Q5XD01
B	-5	VAL	-	expression tag	UNP Q5XD01
B	-4	PRO	-	expression tag	UNP Q5XD01
B	-3	ARG	-	expression tag	UNP Q5XD01
B	-2	GLY	-	expression tag	UNP Q5XD01
B	-1	SER	-	expression tag	UNP Q5XD01
B	0	HIS	-	expression tag	UNP Q5XD01
B	312	ALA	LYS	engineered mutation	UNP Q5XD01
C	-19	MET	-	expression tag	UNP Q5XD01
C	-18	GLY	-	expression tag	UNP Q5XD01
C	-17	SER	-	expression tag	UNP Q5XD01
C	-16	SER	-	expression tag	UNP Q5XD01
C	-15	HIS	-	expression tag	UNP Q5XD01
C	-14	HIS	-	expression tag	UNP Q5XD01
C	-13	HIS	-	expression tag	UNP Q5XD01
C	-12	HIS	-	expression tag	UNP Q5XD01
C	-11	HIS	-	expression tag	UNP Q5XD01
C	-10	HIS	-	expression tag	UNP Q5XD01
C	-9	SER	-	expression tag	UNP Q5XD01
C	-8	SER	-	expression tag	UNP Q5XD01
C	-7	GLY	-	expression tag	UNP Q5XD01
C	-6	LEU	-	expression tag	UNP Q5XD01
C	-5	VAL	-	expression tag	UNP Q5XD01
C	-4	PRO	-	expression tag	UNP Q5XD01
C	-3	ARG	-	expression tag	UNP Q5XD01
C	-2	GLY	-	expression tag	UNP Q5XD01
C	-1	SER	-	expression tag	UNP Q5XD01
C	0	HIS	-	expression tag	UNP Q5XD01
C	312	ALA	LYS	engineered mutation	UNP Q5XD01

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	MET	-	expression tag	UNP Q5XD01
D	-18	GLY	-	expression tag	UNP Q5XD01
D	-17	SER	-	expression tag	UNP Q5XD01
D	-16	SER	-	expression tag	UNP Q5XD01
D	-15	HIS	-	expression tag	UNP Q5XD01
D	-14	HIS	-	expression tag	UNP Q5XD01
D	-13	HIS	-	expression tag	UNP Q5XD01
D	-12	HIS	-	expression tag	UNP Q5XD01
D	-11	HIS	-	expression tag	UNP Q5XD01
D	-10	HIS	-	expression tag	UNP Q5XD01
D	-9	SER	-	expression tag	UNP Q5XD01
D	-8	SER	-	expression tag	UNP Q5XD01
D	-7	GLY	-	expression tag	UNP Q5XD01
D	-6	LEU	-	expression tag	UNP Q5XD01
D	-5	VAL	-	expression tag	UNP Q5XD01
D	-4	PRO	-	expression tag	UNP Q5XD01
D	-3	ARG	-	expression tag	UNP Q5XD01
D	-2	GLY	-	expression tag	UNP Q5XD01
D	-1	SER	-	expression tag	UNP Q5XD01
D	0	HIS	-	expression tag	UNP Q5XD01
D	312	ALA	LYS	engineered mutation	UNP Q5XD01

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

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

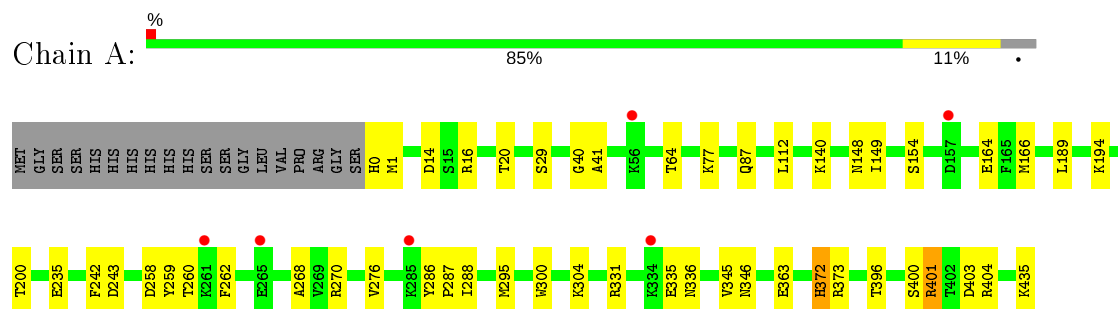
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	165	Total	O	0	0
			165	165		
3	B	127	Total	O	0	0
			127	127		
3	C	175	Total	O	0	0
			175	175		
3	D	110	Total	O	0	0
			110	110		

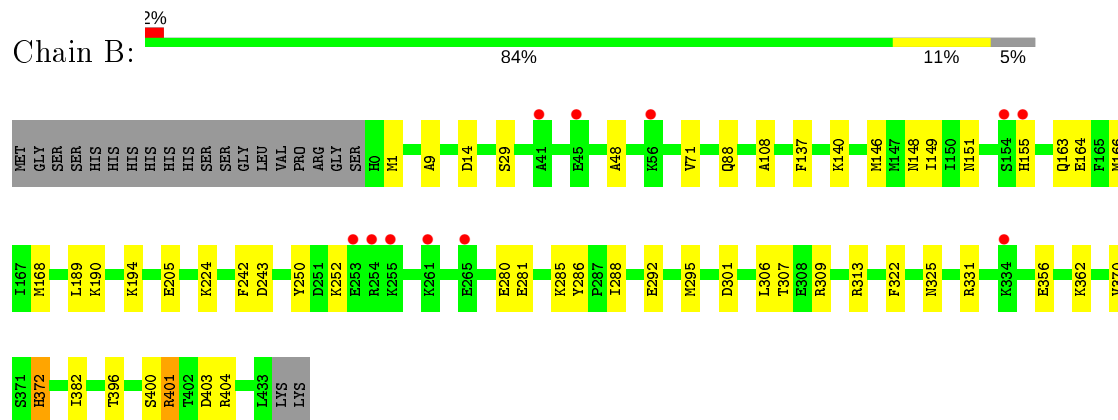
3 Residue-property plots [i](#)

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.

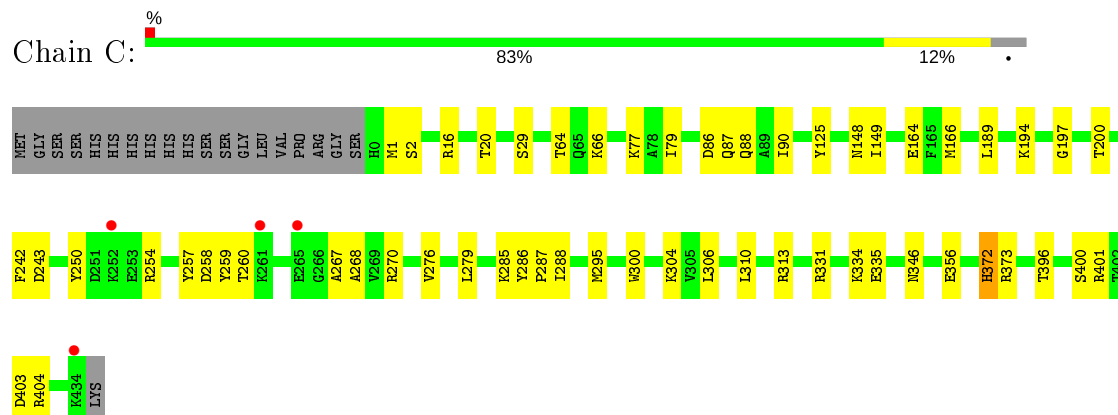
• Molecule 1: ENOLASE



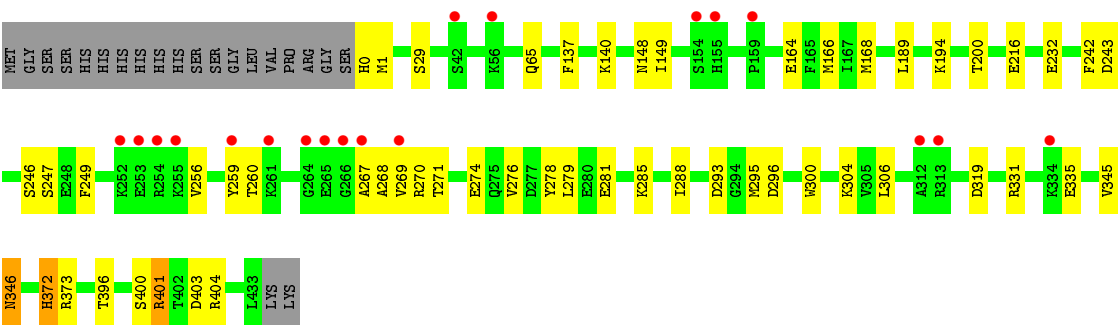
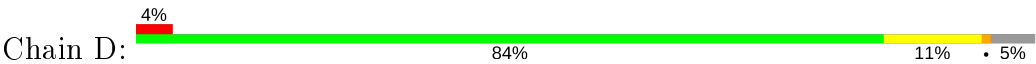
• Molecule 1: ENOLASE



• Molecule 1: ENOLASE



● Molecule 1: ENOLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	181.62Å 181.62Å 56.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	90.81 – 2.15 90.81 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (90.81-2.15) 100.0 (90.81-2.15)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.14Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.172 , 0.199 0.174 , 0.189	Depositor DCC
R_{free} test set	5064 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 13.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.114 for h,-k,-l	Xtriage
Reported twinning fraction	0.140 for h,-k,-l	Depositor
Outliers	0 of 100879 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	27040	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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: PO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.30	0/3392	0.47	0/4585
1	B	0.32	0/3374	0.49	0/4563
1	C	0.31	0/3377	0.47	0/4566
1	D	0.31	0/3374	0.48	0/4563
All	All	0.31	0/13517	0.48	0/18277

There are no bond length outliers.

There are no bond angle outliers.

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	3338	3298	3293	40	0
1	B	3320	3272	3267	37	0
1	C	3324	3279	3275	40	0
1	D	3320	3272	3267	35	0
2	A	10	0	0	2	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
3	A	165	0	0	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	127	0	0	5	0
3	C	175	0	0	7	0
3	D	110	0	0	2	0
All	All	13919	13121	13102	145	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (145) 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:C:295:MET:SD	3:C:2125:HOH:O	2.15	1.03
1:B:301:ASP:OD2	3:B:2097:HOH:O	1.98	0.81
1:A:0:HIS:N	3:A:2001:HOH:O	2.16	0.78
1:D:346:ASN:HD22	1:D:373:ARG:HG3	1.56	0.70
1:C:77:LYS:HG3	3:C:2044:HOH:O	1.90	0.70
1:C:86:ASP:OD2	3:C:2047:HOH:O	2.09	0.69
1:A:0:HIS:CA	3:A:2001:HOH:O	2.41	0.69
1:C:356:GLU:OE2	1:D:140:LYS:NZ	2.26	0.68
1:C:276:VAL:HG22	1:C:295:MET:CE	2.24	0.67
1:A:77:LYS:HG3	3:A:2045:HOH:O	1.94	0.66
1:A:276:VAL:HG22	1:A:295:MET:CE	2.26	0.65
1:B:224:LYS:HG3	3:B:2084:HOH:O	2.00	0.62
1:A:0:HIS:HA	3:A:2001:HOH:O	2.01	0.61
1:B:307:THR:O	3:B:2099:HOH:O	2.16	0.60
1:A:140:LYS:NZ	1:B:356:GLU:OE2	2.36	0.59
1:B:242:PHE:CE2	1:B:288:ILE:HD13	2.39	0.58
1:C:304:LYS:HE2	1:C:335:GLU:HB3	1.87	0.57
1:D:216:GLU:OE2	1:D:285:LYS:NZ	2.34	0.56
1:A:41:ALA:N	2:A:437:PO4:O3	2.39	0.56
1:D:256:VAL:HG11	1:D:269:VAL:HG12	1.88	0.56
1:A:259:TYR:CE2	1:A:270:ARG:NH1	2.74	0.55
1:C:254:ARG:NH2	1:C:258:ASP:OD2	2.39	0.55
1:D:304:LYS:HE2	1:D:335:GLU:HB3	1.87	0.55
1:D:246:SER:HA	1:D:249:PHE:CE2	2.42	0.55
1:D:270:ARG:HB3	1:D:274:GLU:HB2	1.87	0.55
1:B:372:HIS:CD2	1:B:404:ARG:HH11	2.25	0.55
1:D:260:THR:HA	1:D:267:ALA:HB3	1.89	0.54
1:D:346:ASN:ND2	1:D:373:ARG:HG3	2.22	0.54
1:A:304:LYS:HE2	1:A:335:GLU:HB3	1.88	0.54
1:B:285:LYS:HE2	1:B:286:TYR:CZ	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:276:VAL:CG2	1:C:295:MET:CE	2.87	0.52
1:A:140:LYS:HE2	1:B:88:GLN:OE1	2.10	0.52
1:A:154:SER:HA	1:A:262:PHE:CE1	2.45	0.51
1:C:400:SER:O	1:C:401:ARG:HB2	2.10	0.51
1:C:334:LYS:HG3	3:C:2144:HOH:O	2.09	0.51
1:A:346:ASN:OD1	1:A:373:ARG:NH1	2.44	0.51
1:D:279:LEU:HD12	1:D:306:LEU:HD13	1.93	0.51
1:D:1:MET:HE2	1:D:1:MET:HA	1.93	0.51
1:A:276:VAL:CG2	1:A:295:MET:CE	2.89	0.50
1:B:164:GLU:HB2	1:B:243:ASP:HB3	1.92	0.50
1:D:400:SER:O	1:D:401:ARG:HB2	2.12	0.50
1:A:286:TYR:HB3	1:A:287:PRO:HD2	1.92	0.50
1:B:370:VAL:CG1	1:B:382:ILE:HB	2.41	0.50
1:A:1:MET:HB3	1:A:29:SER:HB3	1.94	0.50
1:B:295:MET:HE3	1:B:306:LEU:HD22	1.93	0.50
1:A:336:ASN:HA	3:A:2134:HOH:O	2.12	0.50
1:D:1:MET:HB3	1:D:29:SER:HB3	1.93	0.50
1:D:259:TYR:HB2	1:D:268:ALA:O	2.12	0.50
1:D:256:VAL:HG13	1:D:270:ARG:C	2.33	0.49
1:D:279:LEU:HD12	1:D:306:LEU:CD1	2.43	0.49
1:C:300:TRP:CH2	1:C:331:ARG:HD3	2.48	0.49
1:C:346:ASN:OD1	1:C:373:ARG:NH1	2.44	0.49
1:A:400:SER:O	1:A:401:ARG:HB2	2.13	0.49
1:A:112:LEU:HD11	1:A:345:VAL:CG1	2.43	0.48
1:C:310:LEU:HD23	1:C:313:ARG:NH2	2.28	0.48
1:C:1:MET:HB3	1:C:29:SER:HB3	1.94	0.48
1:A:242:PHE:CE2	1:A:288:ILE:HD13	2.48	0.48
1:C:1:MET:HA	1:C:1:MET:HE2	1.94	0.48
1:A:1:MET:HA	1:A:1:MET:HE2	1.94	0.48
1:A:258:ASP:OD1	1:A:260:THR:OG1	2.26	0.48
1:B:400:SER:O	1:B:401:ARG:HB2	2.14	0.48
1:C:372:HIS:CD2	1:C:404:ARG:HH11	2.32	0.48
1:A:87:GLN:HG2	1:B:137:PHE:CD2	2.49	0.47
1:D:242:PHE:CE2	1:D:288:ILE:HD13	2.50	0.47
1:B:1:MET:HA	1:B:1:MET:HE2	1.97	0.47
1:C:259:TYR:CE2	1:C:270:ARG:NH1	2.83	0.47
1:C:149:ILE:HG12	1:C:189:LEU:CD2	2.45	0.47
1:B:370:VAL:HG23	3:B:2108:HOH:O	2.15	0.47
1:C:242:PHE:CE2	1:C:288:ILE:HD13	2.50	0.47
1:A:372:HIS:CD2	1:A:404:ARG:HH11	2.33	0.46
1:B:1:MET:HB3	1:B:29:SER:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:GLU:OE1	1:B:309:ARG:NH1	2.47	0.46
1:D:300:TRP:CH2	1:D:331:ARG:HD3	2.51	0.46
1:C:16:ARG:NH2	3:C:2008:HOH:O	2.48	0.46
1:B:149:ILE:HG12	1:B:189:LEU:CD2	2.45	0.46
1:B:148:ASN:HA	1:B:166:MET:HG2	1.97	0.46
1:A:112:LEU:HD11	1:A:345:VAL:HG11	1.97	0.46
1:C:88:GLN:OE1	1:D:140:LYS:HE2	2.16	0.46
1:B:148:ASN:ND2	1:B:205:GLU:OE1	2.49	0.46
1:D:278:TYR:O	1:D:281:GLU:HB3	2.16	0.46
1:B:48:ALA:HB1	1:B:108:ALA:HB2	1.97	0.45
1:D:271:THR:OG1	1:D:274:GLU:HG2	2.15	0.45
1:D:372:HIS:CD2	1:D:396:THR:HA	2.52	0.45
1:A:300:TRP:CH2	1:A:331:ARG:HD3	2.51	0.45
1:D:372:HIS:CD2	1:D:404:ARG:HH11	2.35	0.45
1:A:164:GLU:HB2	1:A:243:ASP:HB3	1.99	0.45
1:B:285:LYS:HE2	1:B:286:TYR:CE2	2.52	0.45
1:C:279:LEU:HD12	1:C:306:LEU:HD13	1.99	0.45
1:C:194:LYS:HG3	1:C:200:THR:HG21	1.99	0.45
1:A:148:ASN:HA	1:A:166:MET:HG2	1.99	0.45
1:B:190:LYS:HE2	1:B:194:LYS:HE3	1.98	0.45
1:C:197:GLY:HA2	3:C:2101:HOH:O	2.17	0.45
1:A:16:ARG:NH2	3:A:2013:HOH:O	2.49	0.44
1:B:14:ASP:OD1	1:B:14:ASP:C	2.55	0.44
1:D:194:LYS:HG3	1:D:200:THR:HG21	1.99	0.44
1:A:194:LYS:HG3	1:A:200:THR:HG21	1.99	0.44
1:A:345:VAL:HG23	3:A:2136:HOH:O	2.16	0.44
1:C:260:THR:HA	1:C:267:ALA:HB3	1.98	0.44
1:C:259:TYR:HB2	1:C:268:ALA:O	2.18	0.44
1:C:286:TYR:HB3	1:C:287:PRO:HD2	2.00	0.44
1:B:372:HIS:CD2	1:B:396:THR:C	2.90	0.44
1:A:112:LEU:CD1	1:A:345:VAL:HG12	2.47	0.44
1:A:149:ILE:HG12	1:A:189:LEU:CD2	2.47	0.44
1:B:250:TYR:CE2	1:B:252:LYS:HA	2.53	0.44
1:B:372:HIS:CD2	1:B:396:THR:HA	2.53	0.43
1:B:9:ALA:HB2	1:B:71:VAL:HG11	1.99	0.43
1:D:276:VAL:HG22	1:D:295:MET:CE	2.48	0.43
1:C:148:ASN:HA	1:C:166:MET:HG2	2.00	0.43
1:D:270:ARG:HB3	1:D:274:GLU:CB	2.49	0.43
1:D:148:ASN:HA	1:D:166:MET:HG2	2.00	0.43
1:D:148:ASN:HB2	3:D:2070:HOH:O	2.17	0.43
1:A:372:HIS:CD2	1:A:396:THR:HA	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:GLY:HA3	1:A:373:ARG:NH2	2.33	0.43
1:D:345:VAL:HG23	1:D:346:ASN:H	1.83	0.43
1:D:164:GLU:HB2	1:D:243:ASP:HB3	1.99	0.43
1:C:276:VAL:CG2	1:C:295:MET:HE1	2.49	0.43
1:C:20:THR:HA	1:C:64:THR:HG21	2.01	0.43
1:C:372:HIS:CD2	1:C:396:THR:HA	2.54	0.42
1:D:149:ILE:HG12	1:D:189:LEU:CD2	2.48	0.42
1:B:146:MET:HA	1:B:168:MET:HG2	2.01	0.42
1:A:235:GLU:OE2	1:A:435:LYS:NZ	2.52	0.42
1:C:164:GLU:HB2	1:C:243:ASP:HB3	2.01	0.42
1:D:247:SER:HA	1:D:296:ASP:OD1	2.19	0.42
1:A:14:ASP:C	1:A:14:ASP:OD1	2.58	0.41
1:A:363:GLU:HA	1:B:362:LYS:O	2.20	0.41
1:A:259:TYR:HB2	1:A:268:ALA:O	2.20	0.41
1:B:151:ASN:ND2	3:B:2064:HOH:O	2.50	0.41
1:B:243:ASP:HA	1:B:292:GLU:HB3	2.02	0.41
1:B:155:HIS:CE1	1:B:163:GLN:NE2	2.88	0.41
1:C:276:VAL:HG22	1:C:295:MET:HE3	2.02	0.41
1:D:65:GLN:HG2	3:D:2002:HOH:O	2.20	0.41
1:B:400:SER:O	1:B:401:ARG:CB	2.69	0.41
1:B:331:ARG:NH2	1:C:285:LYS:HD3	2.35	0.41
1:C:87:GLN:HG2	1:D:137:PHE:CD2	2.56	0.41
1:B:322:PHE:O	1:B:325:ASN:HB3	2.20	0.41
1:D:293:ASP:OD2	1:D:319:ASP:HB3	2.21	0.41
1:A:276:VAL:CG2	1:A:295:MET:HE1	2.51	0.40
1:C:2:SER:HB3	1:C:125:TYR:CZ	2.56	0.40
1:C:276:VAL:HG23	1:C:295:MET:HE1	2.03	0.40
2:A:436:PO4:O2	2:A:437:PO4:O4	2.38	0.40
1:A:20:THR:HA	1:A:64:THR:HG21	2.02	0.40
1:C:250:TYR:CD1	1:C:257:TYR:CE1	3.09	0.40
1:C:66:LYS:HE3	3:C:2034:HOH:O	2.19	0.40
1:B:370:VAL:HG11	1:B:382:ILE:HB	2.02	0.40
1:C:79:ILE:HD12	1:C:90:ILE:HG23	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	434/455 (95%)	416 (96%)	17 (4%)	1 (0%)	47	46
1	B	432/455 (95%)	418 (97%)	12 (3%)	2 (0%)	29	22
1	C	433/455 (95%)	417 (96%)	16 (4%)	0	100	100
1	D	432/455 (95%)	414 (96%)	16 (4%)	2 (0%)	29	22
All	All	1731/1820 (95%)	1665 (96%)	61 (4%)	5 (0%)	41	37

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	140	LYS
1	B	401	ARG
1	A	401	ARG
1	D	346	ASN
1	D	401	ARG

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	345/361 (96%)	343 (99%)	2 (1%)	86	90
1	B	343/361 (95%)	339 (99%)	4 (1%)	71	76
1	C	343/361 (95%)	341 (99%)	2 (1%)	86	90
1	D	343/361 (95%)	338 (98%)	5 (2%)	65	69
All	All	1374/1444 (95%)	1361 (99%)	13 (1%)	78	83

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

Mol	Chain	Res	Type
1	A	372	HIS
1	A	403	ASP
1	B	281	GLU
1	B	313	ARG
1	B	372	HIS
1	B	403	ASP
1	C	372	HIS
1	C	403	ASP
1	D	0	HIS
1	D	168	MET
1	D	232	GLU
1	D	372	HIS
1	D	403	ASP

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

Mol	Chain	Res	Type
1	A	163	GLN
1	A	372	HIS
1	B	46	HIS
1	B	155	HIS
1	B	163	GLN
1	B	372	HIS
1	C	138	ASN
1	C	163	GLN
1	C	372	HIS
1	D	65	GLN
1	D	346	ASN
1	D	372	HIS

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 [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	C	437	-	4,4,4	1.06	0	6,6,6	0.50	0
2	PO4	D	437	-	4,4,4	0.96	0	6,6,6	0.33	0
2	PO4	C	436	-	4,4,4	0.78	0	6,6,6	0.56	0
2	PO4	A	436	-	4,4,4	0.82	0	6,6,6	0.46	0
2	PO4	D	436	-	4,4,4	0.81	0	6,6,6	0.42	0
2	PO4	B	436	-	4,4,4	0.90	0	6,6,6	0.36	0
2	PO4	A	437	-	4,4,4	1.10	0	6,6,6	0.33	0
2	PO4	B	437	-	4,4,4	0.96	0	6,6,6	0.48	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	436	PO4	1	0
2	A	437	PO4	2	0

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 ⓘ

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	436/455 (95%)	-0.24	6 (1%) 75 80	17, 31, 59, 88	0
1	B	434/455 (95%)	-0.20	11 (2%) 57 65	17, 31, 62, 97	0
1	C	435/455 (95%)	-0.28	4 (0%) 84 88	17, 28, 52, 76	0
1	D	434/455 (95%)	-0.03	19 (4%) 34 43	17, 32, 78, 116	0
All	All	1739/1820 (95%)	-0.19	40 (2%) 60 68	17, 30, 64, 116	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	267	ALA	5.9
1	D	253	GLU	5.6
1	D	269	VAL	4.9
1	B	155	HIS	4.5
1	D	266	GLY	4.5
1	D	252	LYS	4.0
1	B	154	SER	3.8
1	B	41	ALA	3.6
1	A	157	ASP	3.6
1	A	261	LYS	3.5
1	D	154	SER	3.3
1	D	255	LYS	3.2
1	D	261	LYS	2.9
1	D	312	ALA	2.8
1	A	334	LYS	2.8
1	D	265	GLU	2.8
1	D	155	HIS	2.8
1	D	264	GLY	2.8
1	B	334	LYS	2.7
1	B	261	LYS	2.7
1	D	42	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	253	GLU	2.6
1	D	334	LYS	2.6
1	B	254	ARG	2.6
1	D	254	ARG	2.4
1	D	313	ARG	2.4
1	A	265	GLU	2.4
1	A	56	LYS	2.4
1	A	285	LYS	2.4
1	B	56	LYS	2.4
1	C	265	GLU	2.3
1	C	261	LYS	2.3
1	B	265	GLU	2.3
1	B	45	GLU	2.2
1	C	252	LYS	2.2
1	B	255	LYS	2.2
1	D	159	PRO	2.1
1	D	56	LYS	2.1
1	D	259	TYR	2.1
1	C	434	LYS	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 monosaccharides 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(\AA^2)	Q<0.9
2	PO4	D	436	5/5	0.90	0.18	52,53,63,68	0
2	PO4	A	436	5/5	0.93	0.14	46,49,55,55	0
2	PO4	B	436	5/5	0.96	0.19	44,49,51,56	0
2	PO4	C	436	5/5	0.97	0.09	43,45,48,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PO4	C	437	5/5	0.98	0.15	38,40,45,46	0
2	PO4	D	437	5/5	0.98	0.11	36,40,42,44	0
2	PO4	A	437	5/5	0.98	0.16	40,42,50,52	0
2	PO4	B	437	5/5	0.99	0.07	35,35,37,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.