



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 04:04 PM GMT

PDB ID : 1SP8
Title : 4-HydroxyphenylpyruvateDioxygenase
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Deposited on : 2004-03-16
Resolution : 2.00 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

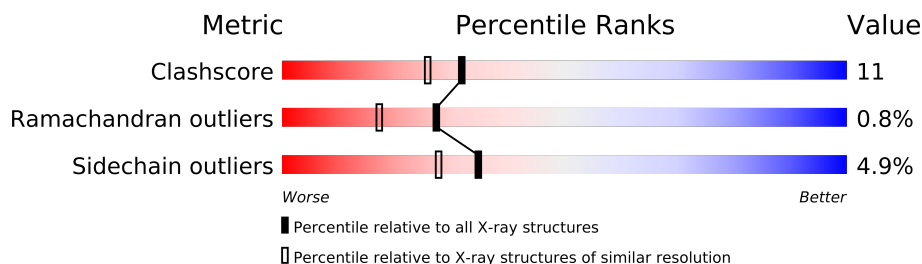
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.00 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))
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	418	
1	B	418	
1	C	418	
1	D	418	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12660 atoms, of which 0 are hydrogen and 0 are deuterium.

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 4-HydroxyphenylpyruvateDioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	137	0	0
			2938	1854	524	550	10			
1	B	392	Total	C	N	O	S	125	0	0
			2971	1876	528	557	10			
1	C	386	Total	C	N	O	S	129	0	0
			2920	1844	519	547	10			
1	D	393	Total	C	N	O	S	130	0	0
			2980	1881	529	560	10			

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	A	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	226	Total	O	0	0
			226	226		
3	B	215	Total	O	0	0
			215	215		
3	C	201	Total	O	0	0
			201	201		
3	D	205	Total	O	1	0
			205	205		

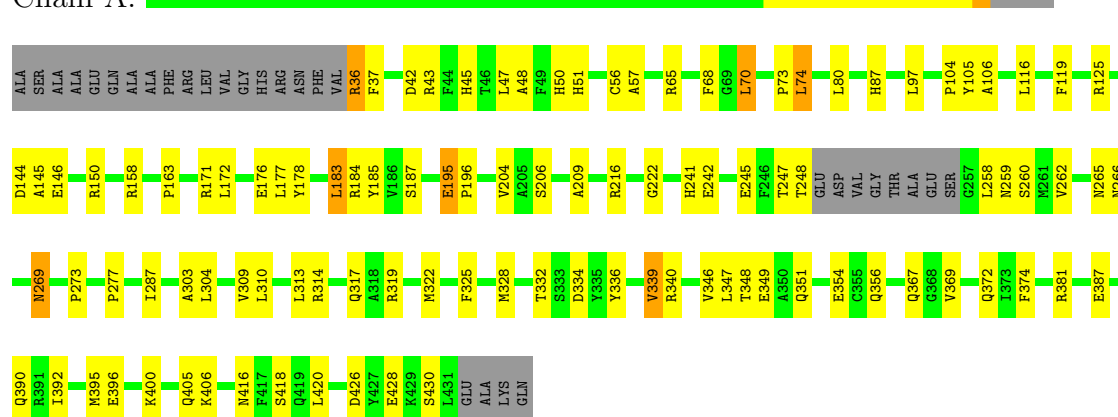
3 Residue-property plots

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.

Note EDS was not executed.

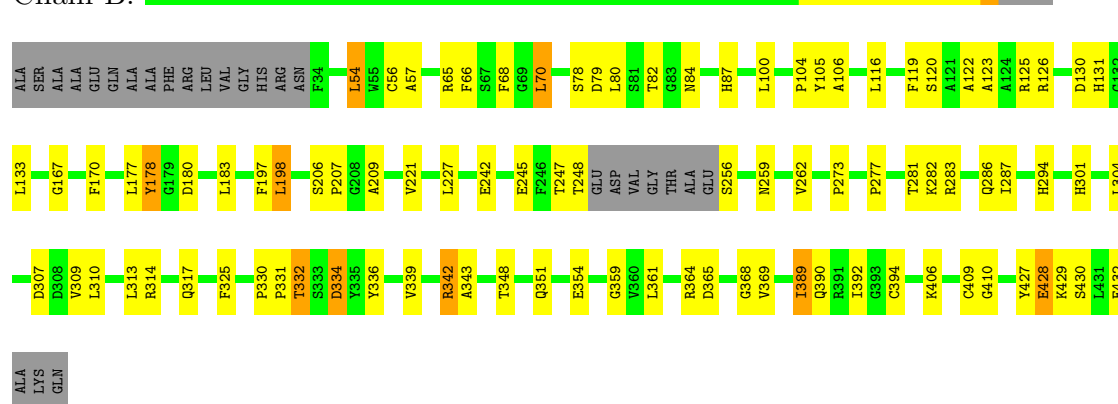
• Molecule 1: 4-HydroxyphenylpyruvateDioxygenase

Chain A:



• Molecule 1: 4-HydroxyphenylpyruvateDioxygenase

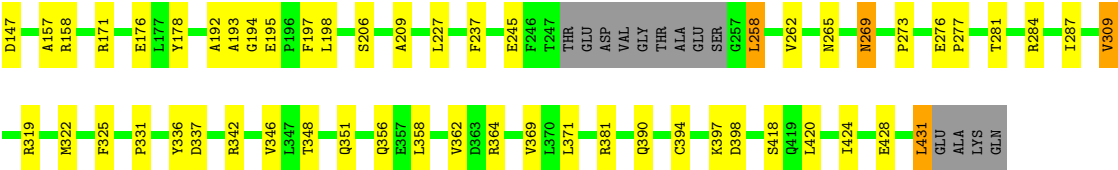
Chain B:



• Molecule 1: 4-HydroxyphenylpyruvateDioxygenase

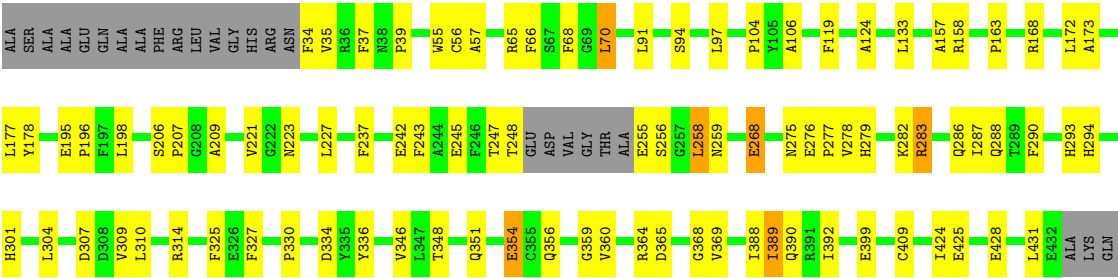
Chain C:





● Molecule 1: 4-HydroxyphenylpyruvateDioxygenase

Chain D:



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.00Å 110.90Å 174.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.84 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (19.84-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.275 , 0.324	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12660	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE2

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	0.52	0/3007	0.72	0/4069
1	B	0.50	0/3041	0.71	1/4115 (0.0%)
1	C	0.51	0/2989	0.70	0/4045
1	D	0.51	1/3050 (0.0%)	0.72	0/4127
All	All	0.51	1/12087 (0.0%)	0.71	1/16356 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	56	CYS	CB-SG	-5.01	1.73	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	54	LEU	CA-CB-CG	5.53	128.02	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2938	0	2830	75	0
1	B	2971	0	2859	70	0
1	C	2920	0	2810	42	0
1	D	2980	0	2865	71	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	226	0	0	4	0
3	B	215	0	0	3	0
3	C	201	0	0	4	0
3	D	205	0	0	8	0
All	All	12660	0	11364	251	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (251) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:206:SER:HB2	1:A:209:ALA:HB3	1.48	0.95
1:D:206:SER:HB2	1:D:209:ALA:HB3	1.48	0.94
1:B:282:LYS:HG3	1:B:283:ARG:H	1.38	0.87
1:C:265:ASN:HD21	1:C:269:ASN:ND2	1.74	0.85
1:C:265:ASN:HD21	1:C:269:ASN:HD21	1.25	0.84
1:B:206:SER:HB2	1:B:209:ALA:HB3	1.59	0.84
1:C:206:SER:HB2	1:C:209:ALA:HB3	1.59	0.84
1:D:195:GLU:HG3	3:D:638:HOH:O	1.81	0.81
1:D:255:GLU:HG3	1:D:256:SER:H	1.46	0.80
1:B:310:LEU:HG	1:B:354:GLU:HG2	1.62	0.80
1:C:262:VAL:HG13	1:C:273:PRO:HG3	1.64	0.79
1:B:79:ASP:H	1:B:82:THR:HG22	1.51	0.76
1:B:332:THR:HG23	1:B:334:ASP:OD2	1.86	0.75
1:A:104:PRO:HB3	1:A:119:PHE:HZ	1.54	0.72
1:A:36:ARG:NH2	1:A:36:ARG:HB2	2.05	0.72
1:D:348:THR:H	1:D:351:GLN:HE21	1.39	0.70
1:B:78:SER:HA	1:B:82:THR:HG21	1.74	0.70
1:A:265:ASN:HD21	1:A:269:ASN:ND2	1.90	0.70
1:B:348:THR:H	1:B:351:GLN:HE21	1.39	0.69
1:C:144:ASP:HB3	1:C:147:ASP:HB3	1.75	0.69
1:A:396:GLU:HB2	1:A:406:LYS:NZ	2.08	0.68
1:A:309:VAL:H	1:A:390:GLN:NE2	1.91	0.68
1:B:339:VAL:HG12	1:B:342:ARG:HH11	1.59	0.68
1:B:365:ASP:HB3	1:B:368:GLY:O	1.93	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:242:GLU:OE1	1:A:245:GLU:HB2	1.94	0.67
1:D:206:SER:HB2	1:D:209:ALA:CB	2.24	0.67
1:A:372:GLN:HE21	1:A:416:ASN:HD22	1.43	0.67
1:B:131:HIS:HE1	1:B:180:ASP:OD2	1.77	0.67
1:B:339:VAL:HA	1:B:342:ARG:HD2	1.77	0.66
1:B:332:THR:HG22	3:B:677:HOH:O	1.95	0.66
1:A:265:ASN:HD21	1:A:269:ASN:HD21	1.43	0.66
1:B:133:LEU:HA	3:B:526:HOH:O	1.94	0.66
1:C:269:ASN:HD22	1:C:269:ASN:H	1.44	0.65
1:D:158:ARG:HH21	1:D:158:ARG:HG3	1.61	0.65
1:A:340:ARG:NH2	1:A:349:GLU:OE1	2.31	0.64
1:C:346:VAL:HG11	1:C:369:VAL:HG21	1.78	0.63
1:A:50:HIS:HD2	1:A:51:HIS:ND1	1.96	0.63
1:B:342:ARG:HG2	1:B:343:ALA:N	2.13	0.63
1:A:51:HIS:HE2	1:A:184:ARG:HH11	1.47	0.62
1:A:336:TYR:HA	1:A:339:VAL:HG13	1.81	0.62
1:C:145:ALA:HB3	1:C:171:ARG:HB3	1.81	0.62
1:D:310:LEU:H	1:D:310:LEU:HD22	1.64	0.62
1:D:258:LEU:HD23	1:D:287:ILE:HD11	1.81	0.62
1:B:79:ASP:O	1:B:82:THR:HG22	1.99	0.62
1:D:248:THR:H	1:D:259:ASN:ND2	1.97	0.61
1:B:369:VAL:HB	1:B:392:ILE:HB	1.83	0.61
1:D:247:THR:HA	1:D:259:ASN:HD22	1.66	0.61
1:A:145:ALA:HB3	1:A:171:ARG:HB3	1.83	0.61
1:B:282:LYS:HG3	1:B:283:ARG:N	2.13	0.60
1:C:106:ALA:HA	1:C:125:ARG:HH11	1.66	0.60
1:A:369:VAL:HB	1:A:392:ILE:HB	1.83	0.60
1:A:247:THR:HA	1:A:259:ASN:HD22	1.65	0.60
1:A:36:ARG:CZ	1:A:36:ARG:HB2	2.32	0.60
1:D:248:THR:H	1:D:259:ASN:HD21	1.51	0.59
1:A:57:ALA:HB3	1:B:57:ALA:HB2	1.85	0.59
1:A:262:VAL:HG13	1:A:273:PRO:HG3	1.85	0.59
1:B:104:PRO:HB3	1:B:119:PHE:HZ	1.68	0.58
1:D:104:PRO:HB3	1:D:119:PHE:HZ	1.68	0.58
1:B:310:LEU:CG	1:B:354:GLU:HG2	2.33	0.58
1:C:342:ARG:NH2	1:C:362:VAL:O	2.36	0.58
1:A:381:ARG:NH1	1:B:106:ALA:HB2	2.19	0.58
1:A:314:ARG:HD3	3:A:633:HOH:O	2.02	0.58
1:D:364:ARG:HG2	1:D:365:ASP:N	2.18	0.58
1:D:248:THR:HG21	1:D:256:SER:OG	2.04	0.57
1:B:348:THR:H	1:B:351:GLN:NE2	2.02	0.57
1:D:172:LEU:HD23	1:D:173:ALA:N	2.19	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:395:MET:SD	1:A:405:GLN:HB2	2.45	0.57
1:C:309:VAL:HB	1:C:390:GLN:HB2	1.87	0.57
1:C:269:ASN:HD22	1:C:269:ASN:N	2.02	0.57
1:A:241:HIS:HD2	3:A:508:HOH:O	1.87	0.57
1:D:283:ARG:NH1	1:D:428:GLU:HB3	2.19	0.57
1:D:327:PHE:HE1	1:D:388:ILE:HD13	1.70	0.56
1:C:104:PRO:HB3	1:C:119:PHE:HZ	1.69	0.56
1:C:192:ALA:C	1:C:194:GLY:H	2.08	0.56
1:A:106:ALA:HA	1:A:125:ARG:HH11	1.71	0.56
1:A:277:PRO:HB3	1:A:287:ILE:HB	1.88	0.55
1:D:348:THR:H	1:D:351:GLN:NE2	2.05	0.55
1:D:277:PRO:HB3	1:D:287:ILE:HB	1.88	0.55
1:A:348:THR:H	1:A:351:GLN:NE2	2.04	0.55
1:A:70:LEU:HD13	1:A:325:PHE:CE1	2.41	0.55
1:A:70:LEU:HD13	1:A:325:PHE:HE1	1.72	0.55
1:C:45:HIS:HB3	1:C:143:ALA:HB2	1.87	0.55
1:B:313:LEU:O	1:B:317:GLN:HG3	2.07	0.55
1:C:122:ALA:HA	3:C:545:HOH:O	2.07	0.55
1:A:73:PRO:HB3	1:A:209:ALA:HB1	1.89	0.55
1:B:131:HIS:HD2	1:B:294:HIS:O	1.90	0.55
1:B:131:HIS:CE1	1:B:180:ASP:OD2	2.59	0.54
1:D:70:LEU:HD13	1:D:325:PHE:CE1	2.42	0.54
1:A:269:ASN:H	1:A:269:ASN:HD22	1.55	0.54
1:A:309:VAL:HB	1:A:390:GLN:HB2	1.88	0.54
1:B:70:LEU:HD13	1:B:325:PHE:CE1	2.42	0.54
1:A:396:GLU:HB2	1:A:406:LYS:HZ2	1.70	0.54
1:D:310:LEU:N	1:D:310:LEU:HD22	2.23	0.54
1:A:309:VAL:H	1:A:390:GLN:HE21	1.56	0.54
1:D:65:ARG:HA	1:D:68:PHE:CE1	2.43	0.54
1:B:177:LEU:HG	1:B:178:TYR:CD2	2.44	0.53
1:B:247:THR:HA	1:B:259:ASN:HD22	1.73	0.53
1:A:158:ARG:HD2	1:A:176:GLU:OE2	2.08	0.53
1:B:336:TYR:O	1:B:339:VAL:HG22	2.08	0.53
1:B:167:GLY:O	1:B:170:PHE:HB2	2.08	0.53
1:B:339:VAL:HG12	1:B:342:ARG:NH1	2.23	0.52
1:C:381:ARG:NH1	1:D:106:ALA:HB2	2.24	0.52
1:D:66:PHE:O	1:D:70:LEU:HB2	2.10	0.52
1:D:330:PRO:HB3	1:D:359:GLY:HA2	1.91	0.52
1:D:221:VAL:HG21	1:D:301:HIS:CE1	2.44	0.52
1:A:269:ASN:N	1:A:269:ASN:HD22	2.07	0.52
1:B:248:THR:HG21	1:B:256:SER:N	2.25	0.52
1:D:307:ASP:HB2	3:D:661:HOH:O	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:394:CYS:HA	3:C:695:HOH:O	2.09	0.52
1:D:346:VAL:HG11	1:D:369:VAL:HG21	1.93	0.51
1:B:332:THR:HG23	1:B:334:ASP:CG	2.30	0.51
1:D:424:ILE:O	1:D:428:GLU:HG2	2.11	0.51
1:B:65:ARG:HA	1:B:68:PHE:CE1	2.46	0.51
1:A:80:LEU:HG	1:A:105:TYR:CZ	2.45	0.51
1:D:365:ASP:OD2	1:D:368:GLY:N	2.42	0.51
1:C:420:LEU:O	1:C:424:ILE:HG23	2.11	0.51
1:D:163:PRO:HA	1:D:172:LEU:O	2.11	0.51
1:D:177:LEU:HG	1:D:178:TYR:CD2	2.46	0.50
1:A:45:HIS:CD2	1:A:47:LEU:HD11	2.47	0.50
1:D:294:HIS:HD2	3:D:513:HOH:O	1.95	0.50
1:C:157:ALA:HB2	1:C:237:PHE:CG	2.47	0.50
1:B:365:ASP:CB	1:B:410:GLY:HA2	2.42	0.50
1:A:178:TYR:CZ	1:A:222:GLY:HA3	2.47	0.49
1:B:294:HIS:HD2	3:B:516:HOH:O	1.95	0.49
1:D:196:PRO:HD2	3:D:637:HOH:O	2.12	0.49
1:A:36:ARG:HG3	1:A:37:PHE:N	2.28	0.49
1:D:258:LEU:HA	1:D:278:VAL:HG23	1.93	0.49
1:A:392:ILE:N	1:A:392:ILE:HD12	2.26	0.49
1:D:35:VAL:O	1:D:35:VAL:HG23	2.13	0.49
1:A:74:LEU:HD13	1:A:204:VAL:HG21	1.94	0.49
1:A:195:GLU:HB2	1:A:196:PRO:HD2	1.95	0.49
1:C:57:ALA:HB3	1:D:57:ALA:HB2	1.93	0.49
1:D:309:VAL:HB	1:D:390:GLN:HB2	1.94	0.49
1:B:330:PRO:HB3	1:B:359:GLY:HA2	1.93	0.49
1:A:322:MET:O	1:B:68:PHE:HB3	2.12	0.49
1:C:336:TYR:CZ	1:C:356:GLN:HA	2.48	0.49
1:A:104:PRO:HB3	1:A:119:PHE:CZ	2.41	0.48
1:B:348:THR:O	1:B:351:GLN:N	2.41	0.48
1:C:81:SER:O	1:D:431:LEU:HD12	2.13	0.48
1:B:314:ARG:HH11	1:B:354:GLU:CD	2.16	0.48
1:B:79:ASP:H	1:B:82:THR:CG2	2.23	0.48
1:B:79:ASP:N	1:B:82:THR:HG22	2.24	0.48
1:A:372:GLN:NE2	1:A:416:ASN:HD22	2.08	0.48
1:C:116:LEU:HD12	3:C:542:HOH:O	2.13	0.48
1:A:183:LEU:HD13	1:A:185:TYR:CE1	2.49	0.48
1:B:364:ARG:HG3	1:B:365:ASP:N	2.29	0.48
1:C:192:ALA:O	1:C:194:GLY:N	2.46	0.48
1:A:258:LEU:N	1:A:258:LEU:HD23	2.28	0.48
1:A:206:SER:HB2	1:A:209:ALA:CB	2.32	0.48
1:B:304:LEU:O	1:B:389:ILE:HD13	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:307:ASP:O	1:B:390:GLN:NE2	2.47	0.48
1:B:262:VAL:HG13	1:B:273:PRO:HG3	1.95	0.48
1:C:227:LEU:HB2	1:C:276:GLU:HB3	1.96	0.48
1:D:286:GLN:OE1	1:D:286:GLN:HA	2.14	0.48
1:A:416:ASN:O	1:A:420:LEU:HD23	2.14	0.47
1:D:133:LEU:HA	3:D:579:HOH:O	2.14	0.47
1:C:192:ALA:C	1:C:194:GLY:N	2.68	0.47
1:C:258:LEU:HG	1:C:258:LEU:O	2.14	0.47
1:D:314:ARG:HH21	1:D:354:GLU:CD	2.16	0.47
1:A:247:THR:O	1:A:247:THR:HG23	2.14	0.47
1:A:97:LEU:HG	1:A:216:ARG:HA	1.95	0.47
1:D:255:GLU:HG3	1:D:256:SER:N	2.23	0.47
1:B:66:PHE:O	1:B:70:LEU:HB2	2.15	0.47
1:D:293:HIS:HD2	3:D:601:HOH:O	1.98	0.47
1:D:255:GLU:HB3	3:D:703:HOH:O	2.14	0.47
1:A:310:LEU:HD13	1:A:354:GLU:HG2	1.97	0.47
1:D:243:PHE:CD2	1:D:268:GLU:HG2	2.51	0.46
1:A:146:GLU:O	1:A:150:ARG:HG3	2.16	0.46
1:B:394:CYS:HB3	1:B:406:LYS:HB3	1.98	0.46
1:C:104:PRO:HB3	1:C:119:PHE:CZ	2.51	0.46
1:C:322:MET:O	1:D:68:PHE:HB3	2.16	0.46
1:B:428:GLU:HA	1:B:428:GLU:OE1	2.15	0.46
1:A:43:ARG:CZ	3:A:597:HOH:O	2.63	0.46
1:D:37:PHE:CD2	1:D:39:PRO:HD3	2.51	0.46
1:B:80:LEU:HG	1:B:105:TYR:CZ	2.51	0.46
1:C:47:LEU:N	1:C:47:LEU:HD12	2.31	0.46
1:B:120:SER:HB3	1:B:123:ALA:HB3	1.98	0.46
1:A:42:ASP:HB2	1:A:266:ASN:ND2	2.31	0.46
1:A:68:PHE:CE1	1:A:319:ARG:HB3	2.51	0.45
1:D:223:ASN:OD1	1:D:275:ASN:HB2	2.16	0.45
1:B:221:VAL:HG21	1:B:301:HIS:CE1	2.51	0.45
1:C:70:LEU:HD13	1:C:325:PHE:CE1	2.52	0.45
1:D:304:LEU:O	1:D:389:ILE:HD13	2.16	0.45
1:A:74:LEU:HD13	1:A:204:VAL:HG11	1.98	0.45
1:D:242:GLU:OE1	1:D:245:GLU:OE1	2.35	0.45
1:D:227:LEU:HB2	1:D:276:GLU:HB3	1.98	0.45
1:B:429:LYS:O	1:B:432:GLU:N	2.49	0.45
1:D:158:ARG:HH21	1:D:158:ARG:CG	2.26	0.45
1:A:177:LEU:HG	1:A:178:TYR:CD2	2.51	0.45
1:D:283:ARG:HH12	1:D:428:GLU:HB3	1.81	0.45
1:D:279:HIS:HE1	1:D:288:GLN:NE2	2.13	0.45
1:D:389:ILE:HD13	1:D:389:ILE:N	2.31	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:389:ILE:HD13	1:D:389:ILE:H	1.82	0.45
1:B:277:PRO:HB3	1:B:287:ILE:HB	1.97	0.45
1:D:94:SER:O	1:D:97:LEU:HB2	2.17	0.44
1:A:346:VAL:HG23	1:A:347:LEU:HG	2.00	0.44
1:B:122:ALA:HA	1:B:125:ARG:NH2	2.32	0.44
1:C:80:LEU:HG	1:C:105:TYR:CZ	2.52	0.44
1:C:431:LEU:N	1:C:431:LEU:HD22	2.31	0.44
1:C:68:PHE:CE1	1:C:319:ARG:HB3	2.52	0.44
1:B:286:GLN:CD	1:B:286:GLN:H	2.20	0.44
1:A:396:GLU:HB2	1:A:406:LYS:HZ1	1.81	0.44
1:A:328:MET:HG3	3:A:592:HOH:O	2.17	0.44
1:D:247:THR:HG23	1:D:259:ASN:ND2	2.32	0.44
1:D:172:LEU:C	1:D:172:LEU:HD23	2.38	0.44
1:D:55:TRP:CE3	1:D:124:ALA:HA	2.53	0.44
1:B:104:PRO:HB3	1:B:119:PHE:CZ	2.52	0.44
1:B:70:LEU:HD13	1:B:325:PHE:HE1	1.80	0.44
1:D:294:HIS:HE1	3:D:519:HOH:O	2.02	0.43
1:C:348:THR:H	1:C:351:GLN:NE2	2.15	0.43
1:A:313:LEU:O	1:A:317:GLN:HG3	2.18	0.43
1:D:157:ALA:HB2	1:D:237:PHE:CG	2.54	0.43
1:D:279:HIS:HE1	1:D:288:GLN:HE21	1.66	0.43
1:A:374:PHE:CE1	1:A:387:GLU:HG3	2.53	0.43
1:A:206:SER:CB	1:A:209:ALA:HB3	2.34	0.43
1:D:158:ARG:NH2	1:D:158:ARG:CG	2.80	0.43
1:C:55:TRP:CE2	1:C:136:ARG:HG2	2.53	0.42
1:B:427:TYR:O	1:B:430:SER:HB3	2.19	0.42
1:A:144:ASP:OD1	1:A:171:ARG:NH2	2.52	0.42
1:B:84:ASN:HD21	1:B:87:HIS:CE1	2.37	0.42
1:C:158:ARG:HD2	1:C:176:GLU:OE2	2.19	0.42
1:A:47:LEU:O	1:A:48:ALA:HB2	2.19	0.42
1:A:245:GLU:HG2	1:A:260:SER:O	2.20	0.42
1:C:77:ARG:HD3	1:C:79:ASP:OD2	2.20	0.42
1:B:283:ARG:HG3	1:B:283:ARG:HH21	1.85	0.41
1:B:309:VAL:HB	1:B:390:GLN:HB2	2.02	0.41
1:B:242:GLU:OE1	1:B:245:GLU:OE1	2.38	0.41
1:C:77:ARG:NH2	3:C:547:HOH:O	2.53	0.41
1:B:126:ARG:O	1:B:130:ASP:HB2	2.21	0.41
1:A:163:PRO:HA	1:A:172:LEU:O	2.21	0.41
1:A:332:THR:OG1	1:A:334:ASP:OD1	2.32	0.41
1:D:168:ARG:HD2	1:D:195:GLU:OE2	2.20	0.41
1:A:245:GLU:HA	1:A:260:SER:O	2.21	0.41
1:D:158:ARG:NH2	1:D:158:ARG:HG3	2.33	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:79:ASP:HB2	1:B:105:TYR:OH	2.21	0.41
1:A:336:TYR:CZ	1:A:356:GLN:HA	2.56	0.41
1:A:171:ARG:HB2	1:A:187:SER:OG	2.21	0.41
1:D:369:VAL:HB	1:D:392:ILE:HB	2.01	0.41
1:D:91:LEU:HD13	1:D:91:LEU:C	2.40	0.41
1:B:100:LEU:HD21	1:B:116:LEU:HD11	2.02	0.41
1:A:348:THR:H	1:A:351:GLN:HE21	1.68	0.41
1:B:197:PHE:CE1	1:B:198:LEU:HD22	2.56	0.41
1:D:290:PHE:C	1:D:290:PHE:CD1	2.94	0.41
1:B:331:PRO:HG2	1:B:361:LEU:CD2	2.51	0.41
1:A:87:HIS:CE1	1:A:116:LEU:HD23	2.56	0.41
1:B:365:ASP:HB2	1:B:410:GLY:HA2	2.03	0.40
1:A:303:ALA:HA	1:A:387:GLU:HB3	2.03	0.40
1:C:331:PRO:HD2	1:C:336:TYR:OH	2.21	0.40
1:B:429:LYS:O	1:B:432:GLU:HA	2.21	0.40
1:D:336:TYR:CZ	1:D:356:GLN:HA	2.57	0.40
1:C:277:PRO:HB3	1:C:287:ILE:HB	2.03	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	384/418 (92%)	362 (94%)	20 (5%)	2 (0%)	38	29
1	B	388/418 (93%)	368 (95%)	18 (5%)	2 (0%)	38	29
1	C	382/418 (91%)	364 (95%)	13 (3%)	5 (1%)	18	8
1	D	389/418 (93%)	365 (94%)	20 (5%)	4 (1%)	22	12
All	All	1543/1672 (92%)	1459 (95%)	71 (5%)	13 (1%)	27	17

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	430	SER

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Mol	Chain	Res	Type
1	C	281	THR
1	C	397	LYS
1	C	193	ALA
1	D	282	LYS
1	C	197	PHE
1	C	398	ASP
1	A	400	LYS
1	B	281	THR
1	D	207	PRO
1	D	399	GLU
1	D	409	CYS
1	B	207	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	292/313 (93%)	277 (95%)	15 (5%)	33	26
1	B	296/313 (95%)	283 (96%)	13 (4%)	39	32
1	C	290/313 (93%)	272 (94%)	18 (6%)	26	18
1	D	297/313 (95%)	286 (96%)	11 (4%)	45	40
All	All	1175/1252 (94%)	1118 (95%)	57 (5%)	35	28

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	ARG
1	A	56	CYS
1	A	65	ARG
1	A	70	LEU
1	A	74	LEU
1	A	183	LEU
1	A	195	GLU
1	A	248	THR
1	A	269	ASN
1	A	304	LEU

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Mol	Chain	Res	Type
1	A	339	VAL
1	A	367	GLN
1	A	418	SER
1	A	426	ASP
1	A	428	GLU
1	B	54	LEU
1	B	56	CYS
1	B	70	LEU
1	B	178	TYR
1	B	183	LEU
1	B	198	LEU
1	B	227	LEU
1	B	332	THR
1	B	334	ASP
1	B	342	ARG
1	B	389	ILE
1	B	409	CYS
1	B	428	GLU
1	C	65	ARG
1	C	70	LEU
1	C	126	ARG
1	C	178	TYR
1	C	195	GLU
1	C	198	LEU
1	C	245	GLU
1	C	258	LEU
1	C	269	ASN
1	C	284	ARG
1	C	309	VAL
1	C	337	ASP
1	C	358	LEU
1	C	364	ARG
1	C	371	LEU
1	C	418	SER
1	C	428	GLU
1	C	431	LEU
1	D	34	PHE
1	D	70	LEU
1	D	198	LEU
1	D	258	LEU
1	D	268	GLU
1	D	283	ARG

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Mol	Chain	Res	Type
1	D	334	ASP
1	D	354	GLU
1	D	360	VAL
1	D	389	ILE
1	D	425	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	HIS
1	A	50	HIS
1	A	241	HIS
1	A	259	ASN
1	A	269	ASN
1	A	351	GLN
1	A	372	GLN
1	A	390	GLN
1	A	405	GLN
1	A	419	GLN
1	B	131	HIS
1	B	259	ASN
1	B	279	HIS
1	B	288	GLN
1	B	294	HIS
1	B	351	GLN
1	B	356	GLN
1	B	419	GLN
1	C	269	ASN
1	C	288	GLN
1	C	351	GLN
1	D	259	ASN
1	D	279	HIS
1	D	288	GLN
1	D	293	HIS
1	D	294	HIS
1	D	351	GLN
1	D	390	GLN
1	D	405	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 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.

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.