



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:14 PM GMT

PDB ID : 1SP8
Title : 4-Hydroxyphenylpyruvate Dioxygenase
Authors : Fritze, I.M.; Linden, L.; Freigang, J.; Auerbach, G.; Huber, R.; Steinbacher, S.
Deposited on : 2004-03-16
Resolution : 2.00 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

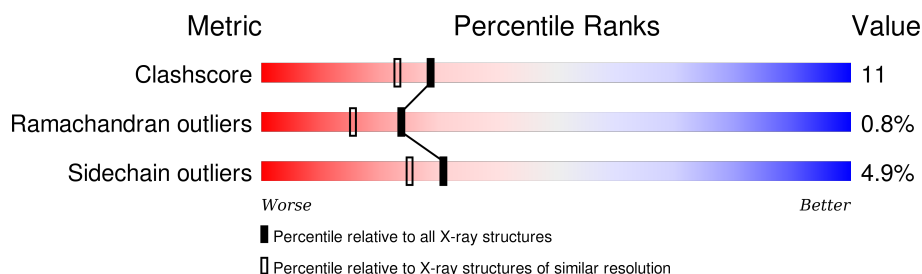
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (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. 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.

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 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 4-Hydroxyphenylpyruvate Dioxygenase.

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.

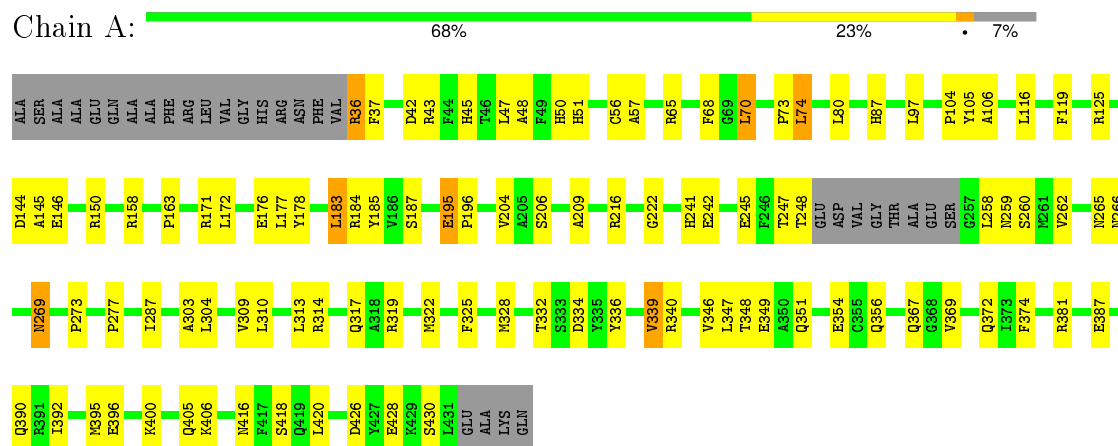
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	226	Total	O	0	0
			226	226		
3	B	215	Total	O	0	0
			215	215		
3	C	201	Total	O	0	0
			201	201		
3	D	205	Total	O	1	0
			205	205		

3 Residue-property plots

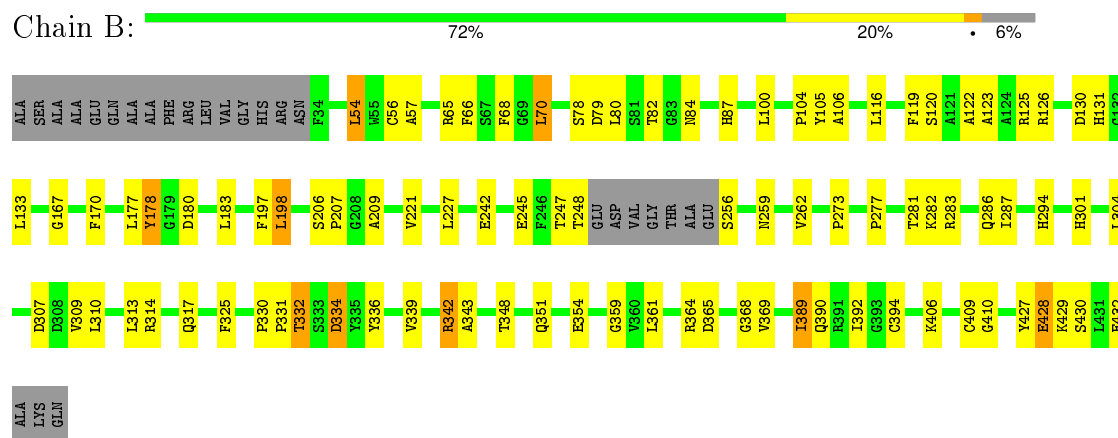
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

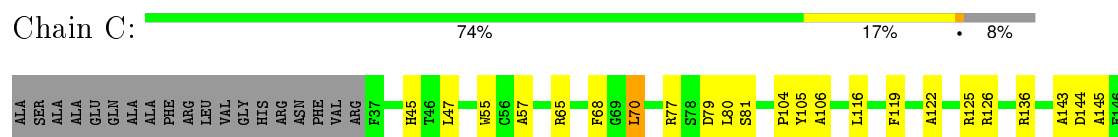
• Molecule 1: 4-Hydroxyphenylpyruvate Dioxygenase

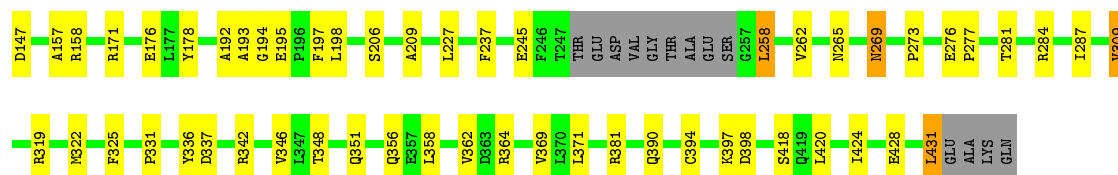


• Molecule 1: 4-Hydroxyphenylpyruvate Dioxygenase



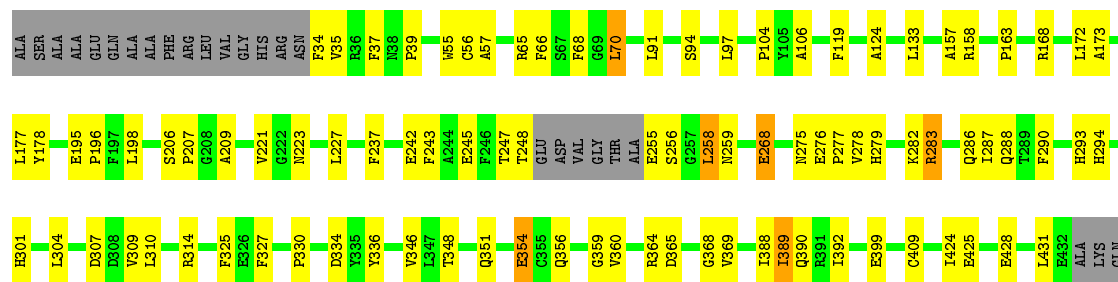
• Molecule 1: 4-Hydroxyphenylpyruvate Dioxygenase





• Molecule 1: 4-Hydroxyphenylpyruvate Dioxygenase

Chain D: 72% 21% 6%



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	89.00 Å 110.90 Å 174.80 Å 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 [i](#)

5.1 Standard geometry [i](#)

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 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	2938	0	2830	75	0
1	B	2971	0	2859	70	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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

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 11.

All (251) 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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:365:ASP:HB3	1:B:368:GLY:O	1.93	0.67
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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 [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	384/418 (92%)	362 (94%)	20 (5%)	2 (0%)	34 26
1	B	388/418 (93%)	368 (95%)	18 (5%)	2 (0%)	34 26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	382/418 (91%)	364 (95%)	13 (3%)	5 (1%)	15	7
1	D	389/418 (93%)	365 (94%)	20 (5%)	4 (1%)	19	11
All	All	1543/1672 (92%)	1459 (95%)	71 (5%)	13 (1%)	24	15

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	430	SER
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%)	29	23
1	B	296/313 (95%)	283 (96%)	13 (4%)	35	30
1	C	290/313 (93%)	272 (94%)	18 (6%)	23	16
1	D	297/313 (95%)	286 (96%)	11 (4%)	41	38
All	All	1175/1252 (94%)	1118 (95%)	57 (5%)	31	25

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
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

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Mol	Chain	Res	Type
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
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

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Mol	Chain	Res	Type
1	D	293	HIS
1	D	294	HIS
1	D	351	GLN
1	D	390	GLN
1	D	405	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates 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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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