



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:42 PM UTC

PDB ID : 3WRC / pdb_00003wrc
Title : Crystal structure of the anaerobic DesB-Protocatechuate (PCA) complex
Authors : Sugimoto, K.; Senda, M.; Kasai, D.; Fukuda, M.; Masai, E.; Senda, T.
Deposited on : 2014-02-21
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

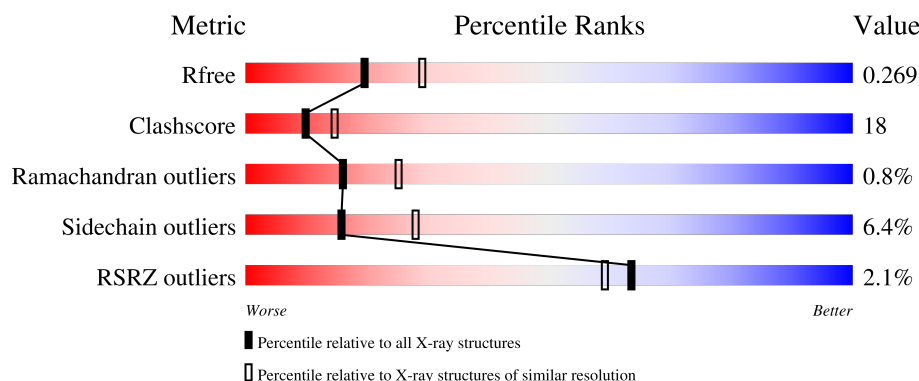
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

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

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6433 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.

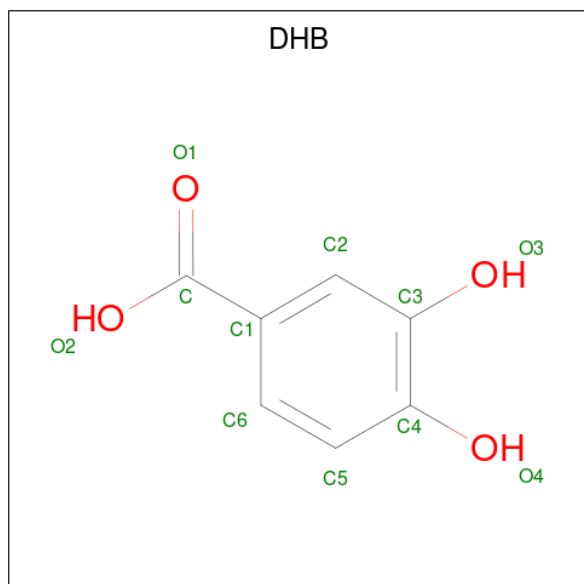
- Molecule 1 is a protein called Gallate dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	0	0	0
			3169	2033	545	577	14			
1	B	404	Total	C	N	O	S	0	0	0
			3209	2062	549	584	14			

- Molecule 2 is FE (II) ION (CCD ID: FE2) (formula: Fe).

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

- Molecule 3 is 3,4-DIHYDROXYBENZOIC ACID (CCD ID: DHB) (formula: C₇H₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			11	7	4		

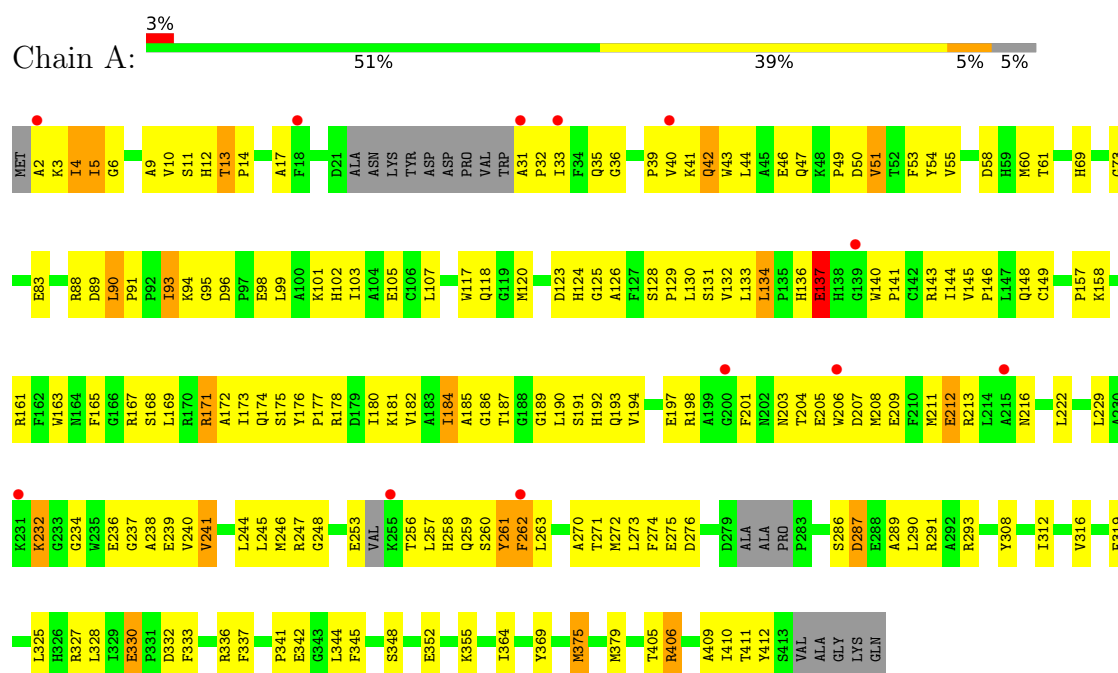
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	10	Total	O	0	0
			10	10		
4	B	32	Total	O	0	0
			32	32		

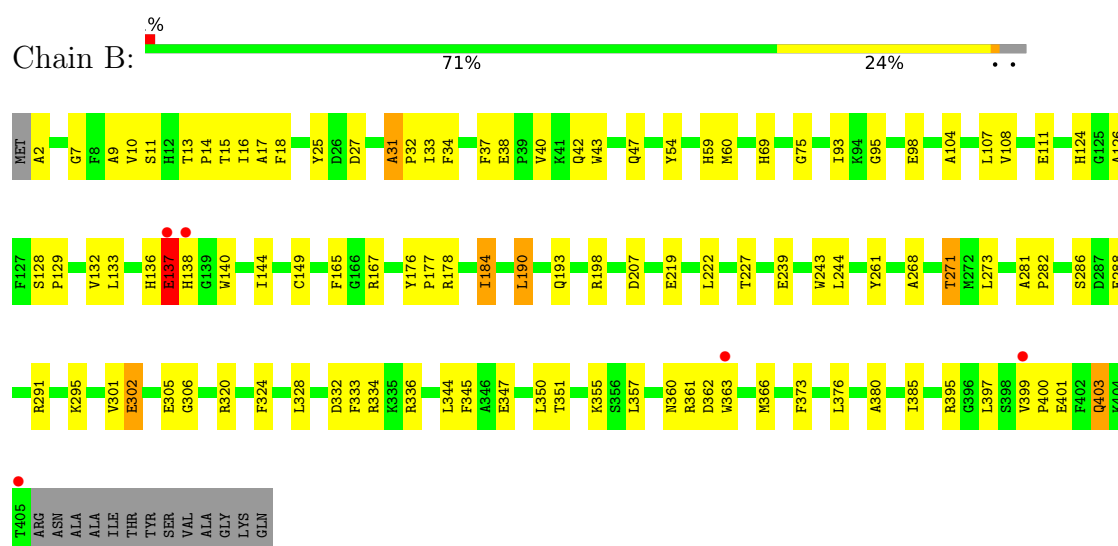
3 Residue-property plots [i](#)

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: Gallate dioxygenase



• Molecule 1: Gallate dioxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.83Å 61.01Å 117.55Å 90.00° 97.84° 90.00°	Depositor
Resolution (Å)	54.04 – 2.40 54.04 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (54.04-2.40) 97.9 (54.04-2.40)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.94 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.4.0078	Depositor
R, R_{free}	0.217 , 0.276 (Not available) , 0.269	Depositor DCC
R_{free} test set	1568 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.365	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6433	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% 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

5.1 Standard geometry

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

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	0.57	1/3257 (0.0%)	0.87	2/4408 (0.0%)
1	B	0.71	0/3304	0.92	4/4482 (0.1%)
All	All	0.64	1/6561 (0.0%)	0.89	6/8890 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	232	LYS	CG-CD	6.25	1.71	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	330	GLU	CA-C-N	7.69	128.08	119.32
1	A	330	GLU	C-N-CA	7.69	128.08	119.32
1	B	243	TRP	N-CA-C	-5.23	105.76	111.82
1	B	59	HIS	N-CA-C	5.03	116.81	109.71
1	B	31	ALA	CA-C-N	5.03	124.81	119.28
1	B	31	ALA	C-N-CA	5.03	124.81	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3169	0	3059	158	0
1	B	3209	0	3093	65	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	B	11	0	3	0	0
4	A	10	0	0	0	0
4	B	32	0	0	2	0
All	All	6433	0	6155	220	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 18.

All (220) 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:4:ILE:HD13	1:A:4:ILE:H	1.13	1.05
1:A:33:ILE:HD11	1:A:262:PHE:HB2	1.45	0.94
1:A:167:ARG:NH1	1:A:248:GLY:O	2.05	0.90
1:B:13:THR:HG22	1:B:15:THR:H	1.37	0.88
1:A:4:ILE:HD13	1:A:4:ILE:N	1.89	0.87
1:A:175:SER:HB3	1:A:290:LEU:HD23	1.57	0.86
1:A:241:VAL:HA	1:A:244:LEU:HD13	1.60	0.84
1:A:90:LEU:HD13	1:A:91:PRO:HD2	1.59	0.83
1:A:44:LEU:HD13	1:A:134:LEU:HD23	1.62	0.82
1:A:4:ILE:H	1:A:4:ILE:CD1	1.95	0.79
1:A:39:PRO:HG2	1:A:258:HIS:CG	2.18	0.79
1:B:98:GLU:OE2	1:B:178:ARG:NH1	2.16	0.77
1:A:39:PRO:HG2	1:A:258:HIS:ND1	2.00	0.77
1:A:352:GLU:HA	1:A:355:LYS:HG3	1.68	0.75
1:A:134:LEU:HB3	1:A:141:PRO:HG3	1.66	0.75
1:A:36:GLY:O	1:A:258:HIS:NE2	2.20	0.75
1:A:406:ARG:HD2	1:A:406:ARG:N	2.04	0.73
1:B:128:SER:HB3	1:B:129:PRO:HD3	1.70	0.73
1:B:302:GLU:CD	1:B:302:GLU:H	1.95	0.72
1:B:37:PHE:CD2	1:B:132:VAL:HG11	2.24	0.72
1:A:88:ARG:HH12	1:A:128:SER:HA	1.53	0.72
1:A:128:SER:HB2	1:A:129:PRO:HD3	1.69	0.72
1:A:260:SER:O	1:A:261:TYR:HB2	1.88	0.71
1:A:32:PRO:HA	1:A:35:GLN:HB2	1.72	0.71
1:A:257:LEU:HD11	1:A:275:GLU:HG3	1.74	0.70
1:A:17:ALA:HB2	1:A:128:SER:OG	1.92	0.68
1:A:96:ASP:OD2	1:A:99:LEU:HB2	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:HIS:O	1:A:137:GLU:C	2.38	0.67
1:A:2:ALA:HB2	1:A:178:ARG:O	1.94	0.66
1:A:327:ARG:NH2	1:A:330:GLU:OE1	2.24	0.66
1:A:98:GLU:OE1	1:A:178:ARG:NH2	2.30	0.64
1:A:163:TRP:HZ3	1:A:247:ARG:HG3	1.62	0.64
1:B:403:GLN:NE2	1:B:403:GLN:O	2.29	0.64
1:A:12:HIS:HB2	1:A:125:GLY:HA2	1.79	0.64
1:A:51:VAL:HG23	1:A:182:VAL:HG13	1.79	0.64
1:A:345:PHE:HB3	1:A:355:LYS:HD3	1.79	0.64
1:A:175:SER:O	1:A:291:ARG:NH1	2.32	0.63
1:B:10:VAL:HG22	1:B:11:SER:O	1.99	0.62
1:A:193:GLN:HA	1:A:236:GLU:OE1	1.99	0.62
1:A:10:VAL:HG12	1:A:187:THR:HG22	1.82	0.61
1:A:171:ARG:NH1	1:A:287:ASP:OD1	2.31	0.60
1:A:4:ILE:N	1:A:4:ILE:CD1	2.56	0.60
1:A:203:ASN:HB3	1:A:206:TRP:HB3	1.84	0.60
1:A:51:VAL:CG2	1:A:182:VAL:HG22	2.31	0.60
1:A:69:HIS:CD2	1:A:157:PRO:HD3	2.37	0.60
1:A:171:ARG:HH22	1:A:287:ASP:CG	2.10	0.59
1:B:13:THR:HG22	1:B:15:THR:N	2.14	0.59
1:B:31:ALA:N	1:B:32:PRO:HD2	2.17	0.59
1:A:51:VAL:HG13	1:A:143:ARG:NH1	2.18	0.58
1:A:41:LYS:HA	1:A:44:LEU:HD12	1.86	0.58
1:A:172:ALA:O	1:A:175:SER:OG	2.20	0.58
1:A:51:VAL:HG23	1:A:182:VAL:HG22	1.86	0.57
1:A:333:PHE:CE1	1:A:344:LEU:HD21	2.39	0.57
1:B:336:ARG:HD3	1:B:344:LEU:HD13	1.86	0.57
1:A:5:ILE:HD11	1:A:49:PRO:HB3	1.87	0.57
1:A:39:PRO:HG2	1:A:258:HIS:CE1	2.39	0.57
1:A:177:PRO:O	1:A:178:ARG:HG2	2.05	0.56
1:A:44:LEU:HD13	1:A:134:LEU:CD2	2.32	0.56
1:A:241:VAL:HA	1:A:244:LEU:CD1	2.35	0.56
1:A:4:ILE:HG22	1:A:182:VAL:HB	1.87	0.56
1:A:229:LEU:HB3	1:A:238:ALA:HA	1.87	0.56
1:A:43:TRP:CD1	1:A:47:GLN:HE21	2.24	0.55
1:B:328:LEU:O	1:B:334:ARG:HD3	2.06	0.55
1:B:9:ALA:O	1:B:271:THR:HA	2.06	0.55
1:A:55:VAL:HB	1:A:186:GLY:HA2	1.88	0.55
1:B:2:ALA:HB2	1:B:176:TYR:O	2.07	0.55
1:A:189:GLY:HA2	1:A:192:HIS:HE1	1.71	0.55
1:B:261:TYR:CE1	1:B:268:ALA:HB1	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:TYR:CE2	1:A:126:ALA:HA	2.42	0.54
1:A:118:GLN:HB2	1:A:319:PHE:CE1	2.43	0.54
1:B:397:LEU:HD22	1:B:401:GLU:HB3	1.90	0.54
1:A:128:SER:CB	1:A:129:PRO:HD3	2.39	0.53
1:A:189:GLY:HA2	1:A:192:HIS:CE1	2.44	0.53
1:A:257:LEU:HD11	1:A:275:GLU:CG	2.39	0.53
1:B:288:GLU:N	1:B:288:GLU:OE2	2.42	0.53
1:A:90:LEU:HD13	1:A:91:PRO:CD	2.37	0.52
1:A:236:GLU:O	1:A:239:GLU:HG2	2.10	0.52
1:A:325:LEU:HB2	1:A:375:MET:HG3	1.90	0.52
1:B:104:ALA:O	1:B:108:VAL:HG23	2.10	0.52
1:A:43:TRP:CE2	1:A:47:GLN:HG3	2.45	0.52
1:B:362:ASP:O	1:B:366:MET:HG3	2.10	0.52
1:A:54:TYR:HE2	1:A:126:ALA:HA	1.75	0.52
1:A:3:LYS:O	1:A:181:LYS:HA	2.10	0.51
1:A:54:TYR:HA	1:A:185:ALA:HB3	1.92	0.51
1:A:209:GLU:OE2	1:A:213:ARG:HD2	2.10	0.51
1:A:312:ILE:O	1:A:316:VAL:HG23	2.09	0.51
1:B:357:LEU:O	1:B:366:MET:HE2	2.11	0.51
1:A:337:PHE:O	1:A:341:PRO:HG3	2.09	0.51
1:A:53:PHE:HE1	1:A:103:ILE:HD13	1.75	0.51
1:A:95:GLY:HA2	1:A:144:ILE:O	2.10	0.50
1:A:163:TRP:CZ3	1:A:247:ARG:HG3	2.45	0.50
1:A:342:GLU:OE1	1:A:342:GLU:HA	2.11	0.50
1:A:51:VAL:CG1	1:A:143:ARG:NH1	2.74	0.50
1:B:128:SER:HB3	1:B:129:PRO:CD	2.39	0.50
1:A:163:TRP:CD1	1:A:222:LEU:HD11	2.47	0.50
1:A:236:GLU:OE1	1:A:236:GLU:N	2.42	0.50
1:A:257:LEU:HD11	1:A:275:GLU:CD	2.37	0.50
1:B:17:ALA:HB2	1:B:128:SER:HA	1.94	0.49
1:A:171:ARG:HH21	1:A:174:GLN:HB3	1.77	0.49
1:A:237:GLY:O	1:A:240:VAL:HG22	2.11	0.49
1:B:43:TRP:CE2	1:B:47:GLN:HG3	2.48	0.49
1:A:90:LEU:HD11	1:A:140:TRP:CZ2	2.48	0.49
1:A:163:TRP:O	1:A:167:ARG:HG2	2.13	0.48
1:A:175:SER:HB3	1:A:290:LEU:CD2	2.38	0.48
1:B:13:THR:CG2	1:B:15:THR:H	2.20	0.48
1:A:134:LEU:HB3	1:A:141:PRO:CG	2.38	0.48
1:A:190:LEU:HD22	1:A:207:ASP:CG	2.38	0.48
1:B:93:ILE:HB	1:B:140:TRP:CD1	2.49	0.48
1:A:102:HIS:NE2	1:A:175:SER:OG	2.33	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:HIS:O	1:A:259:GLN:HG2	2.14	0.48
1:B:7:GLY:HA2	1:B:184:ILE:O	2.14	0.47
1:B:43:TRP:CD2	1:B:273:LEU:HD11	2.49	0.47
1:A:9:ALA:HB2	1:A:246:MET:HE3	1.96	0.47
1:B:320:ARG:HH21	1:B:351:THR:HG23	1.78	0.47
1:B:190:LEU:HD13	1:B:207:ASP:HB3	1.97	0.47
1:B:345:PHE:HB3	1:B:355:LYS:HD3	1.96	0.47
1:A:332:ASP:O	1:A:336:ARG:HG3	2.15	0.47
1:A:10:VAL:HG22	1:A:11:SER:N	2.30	0.47
1:B:37:PHE:HD2	1:B:132:VAL:HG11	1.74	0.47
1:B:40:VAL:HG12	1:B:133:LEU:HD22	1.96	0.47
1:A:204:THR:HA	1:A:207:ASP:HB2	1.96	0.47
1:B:18:PHE:CD1	1:B:18:PHE:C	2.92	0.47
1:B:75:GLY:HA3	4:B:627:HOH:O	2.14	0.46
1:A:197:GLU:OE1	1:B:306:GLY:HA2	2.15	0.46
1:A:286:SER:O	1:A:289:ALA:N	2.48	0.46
1:B:37:PHE:CD1	1:B:271:THR:HG21	2.51	0.46
1:A:208:MET:O	1:A:212:GLU:HB3	2.15	0.46
1:A:90:LEU:HD11	1:A:140:TRP:HZ2	1.80	0.46
1:A:93:ILE:HG12	1:A:94:LYS:N	2.31	0.46
1:B:136:HIS:O	1:B:138:HIS:N	2.48	0.46
1:B:31:ALA:HB3	1:B:32:PRO:HD3	1.97	0.46
1:A:73:GLY:HA3	1:A:146:PRO:HG2	1.98	0.46
1:B:167:ARG:HD2	1:B:167:ARG:HA	1.61	0.46
1:B:244:LEU:HA	1:B:244:LEU:HD23	1.75	0.46
1:A:134:LEU:HB2	1:A:140:TRP:CZ3	2.51	0.45
1:B:14:PRO:HD3	1:B:124:HIS:CD2	2.51	0.45
1:B:40:VAL:HG22	1:B:273:LEU:HB2	1.96	0.45
1:B:328:LEU:HA	1:B:333:PHE:HD2	1.81	0.45
1:A:405:THR:C	1:A:406:ARG:HD2	2.41	0.45
1:B:54:TYR:CE2	1:B:126:ALA:HA	2.52	0.45
1:A:130:LEU:HA	1:A:133:LEU:HB2	1.97	0.45
1:A:369:TYR:O	1:A:369:TYR:CD2	2.70	0.45
1:A:128:SER:O	1:A:131:SER:OG	2.34	0.45
1:A:31:ALA:C	1:A:33:ILE:H	2.24	0.45
1:A:94:LYS:O	1:A:144:ILE:HG22	2.16	0.45
1:A:261:TYR:CD2	1:A:262:PHE:N	2.84	0.45
1:A:42:GLN:HG3	1:A:43:TRP:N	2.31	0.45
1:A:96:ASP:OD2	1:A:99:LEU:CB	2.64	0.45
1:A:163:TRP:HZ3	1:A:247:ARG:CG	2.29	0.45
1:A:257:LEU:O	1:A:258:HIS:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:MET:HE1	1:B:60:MET:HB3	1.97	0.45
1:B:399:VAL:HB	1:B:400:PRO:HD3	1.98	0.45
1:A:409:ALA:C	1:A:410:ILE:HG23	2.42	0.45
1:A:333:PHE:CZ	1:A:344:LEU:HD21	2.53	0.44
1:A:201:PHE:HE2	1:A:232:LYS:O	1.99	0.44
1:A:47:GLN:H	1:A:47:GLN:HG2	1.63	0.44
1:A:90:LEU:HA	1:A:90:LEU:HD22	1.65	0.44
1:A:184:ILE:HG22	1:A:184:ILE:O	2.16	0.44
1:B:177:PRO:HG3	1:B:291:ARG:HD2	1.99	0.44
1:A:161:ARG:HD3	1:B:111:GLU:HG2	2.00	0.44
1:A:364:ILE:HD12	1:A:364:ILE:HA	1.86	0.44
1:B:38:GLU:O	1:B:42:GLN:HG3	2.18	0.44
1:A:107:LEU:HD13	1:A:165:PHE:CE1	2.53	0.44
1:A:101:LYS:HE2	1:A:105:GLU:OE2	2.18	0.43
1:A:176:TYR:CE2	1:A:178:ARG:HG3	2.53	0.43
1:B:25:TYR:HA	1:B:34:PHE:CD2	2.54	0.43
1:A:328:LEU:HD13	1:A:379:MET:HA	2.00	0.43
1:B:95:GLY:HA2	1:B:144:ILE:O	2.19	0.43
1:B:107:LEU:HD13	1:B:165:PHE:CZ	2.53	0.43
1:A:6:GLY:N	1:A:276:ASP:HB2	2.34	0.43
1:B:136:HIS:C	1:B:138:HIS:H	2.26	0.43
1:B:295:LYS:HA	1:B:295:LYS:HD3	1.76	0.43
1:A:13:THR:HG22	1:A:14:PRO:HD2	2.00	0.43
1:A:43:TRP:CB	1:A:273:LEU:HD21	2.48	0.43
1:B:363:TRP:CH2	1:B:380:ALA:HB2	2.54	0.43
1:A:165:PHE:CD2	1:A:245:LEU:HD11	2.54	0.42
1:A:198:ARG:NH1	1:A:234:GLY:HA2	2.33	0.42
1:A:50:ASP:HB2	1:A:181:LYS:O	2.19	0.42
1:B:219:GLU:HA	1:B:222:LEU:HG	2.00	0.42
1:A:99:LEU:HG	1:A:103:ILE:HD11	2.01	0.42
1:B:239:GLU:HB3	4:B:604:HOH:O	2.18	0.42
1:A:69:HIS:O	1:A:149:CYS:HA	2.19	0.42
1:A:176:TYR:HA	1:A:177:PRO:HD2	1.92	0.42
1:A:211:MET:CE	1:A:270:ALA:HB1	2.48	0.42
1:A:107:LEU:HD23	1:A:168:SER:HB3	2.00	0.42
1:A:260:SER:HB2	1:A:271:THR:OG1	2.18	0.42
1:B:136:HIS:C	1:B:138:HIS:N	2.78	0.42
1:A:69:HIS:NE2	1:A:157:PRO:HD3	2.35	0.42
1:A:163:TRP:CD1	1:A:222:LEU:CD1	3.03	0.42
1:B:137:GLU:OE1	1:B:137:GLU:HA	2.20	0.42
1:A:411:THR:O	1:A:412:TYR:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:HIS:O	1:B:149:CYS:HA	2.19	0.42
1:A:169:LEU:O	1:A:173:ILE:HG13	2.19	0.41
1:A:216:ASN:N	1:A:216:ASN:ND2	2.64	0.41
1:B:324:PHE:CG	1:B:350:LEU:HD11	2.55	0.41
1:A:117:TRP:HB2	1:A:120:MET:CE	2.50	0.41
1:A:5:ILE:HG23	1:A:181:LYS:HB3	2.01	0.41
1:B:193:GLN:OE1	1:B:198:ARG:HD3	2.20	0.41
1:A:51:VAL:HG13	1:A:143:ARG:HH12	1.84	0.41
1:A:189:GLY:HA3	1:A:190:LEU:HA	1.75	0.41
1:A:244:LEU:HD12	1:A:244:LEU:H	1.85	0.41
1:A:257:LEU:CD1	1:A:275:GLU:HG3	2.46	0.41
1:A:186:GLY:HA3	1:A:246:MET:HB2	2.03	0.41
1:A:287:ASP:OD1	1:A:287:ASP:N	2.52	0.41
1:A:308:TYR:CD2	1:A:308:TYR:N	2.89	0.41
1:A:144:ILE:HG12	1:A:145:VAL:N	2.36	0.41
1:A:40:VAL:HG22	1:A:273:LEU:HD13	2.02	0.40
1:A:158:LYS:O	1:A:161:ARG:HB3	2.21	0.40
1:B:360:ASN:O	1:B:361:ARG:C	2.65	0.40
1:A:83:GLU:CD	1:A:123:ASP:HB2	2.46	0.40
1:A:117:TRP:CE3	1:A:120:MET:HE3	2.57	0.40
1:A:191:SER:O	1:A:192:HIS:HB3	2.21	0.40
1:B:373:PHE:O	1:B:376:LEU:N	2.54	0.40
1:A:14:PRO:HB3	1:A:124:HIS:CG	2.55	0.40
1:A:58:ASP:H	1:A:148:GLN:CD	2.29	0.40
1:A:272:MET:HB3	1:A:274:PHE:CE1	2.56	0.40
1:B:33:ILE:H	1:B:33:ILE:HG12	1.74	0.40
1:A:5:ILE:O	1:A:43:TRP:HZ2	2.04	0.40
1:A:194:VAL:HG22	1:A:236:GLU:OE2	2.21	0.40
1:B:281:ALA:HB1	1:B:282:PRO:HD2	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	391/418 (94%)	356 (91%)	31 (8%)	4 (1%)	12	20
1	B	402/418 (96%)	384 (96%)	16 (4%)	2 (0%)	24	37
All	All	793/836 (95%)	740 (93%)	47 (6%)	6 (1%)	16	25

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	GLU
1	A	261	TYR
1	B	190	LEU
1	A	46	GLU
1	A	89	ASP
1	B	137	GLU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	326/340 (96%)	299 (92%)	27 (8%)	10	17
1	B	330/340 (97%)	315 (96%)	15 (4%)	24	42
All	All	656/680 (96%)	614 (94%)	42 (6%)	16	28

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	A	5	ILE
1	A	13	THR
1	A	42	GLN
1	A	51	VAL
1	A	60	MET
1	A	61	THR
1	A	90	LEU
1	A	93	ILE
1	A	132	VAL

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Mol	Chain	Res	Type
1	A	134	LEU
1	A	137	GLU
1	A	171	ARG
1	A	180	ILE
1	A	184	ILE
1	A	205	GLU
1	A	212	GLU
1	A	241	VAL
1	A	253	GLU
1	A	256	THR
1	A	262	PHE
1	A	263	LEU
1	A	287	ASP
1	A	293	ARG
1	A	348	SER
1	A	375	MET
1	A	406	ARG
1	B	16	ILE
1	B	27	ASP
1	B	137	GLU
1	B	184	ILE
1	B	227	THR
1	B	271	THR
1	B	286	SER
1	B	301	VAL
1	B	302	GLU
1	B	305	GLU
1	B	332	ASP
1	B	347	GLU
1	B	385	ILE
1	B	395	ARG
1	B	403	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	164	ASN
1	A	192	HIS
1	A	202	ASN
1	A	216	ASN
1	A	360	ASN

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Mol	Chain	Res	Type
1	B	23	ASN
1	B	216	ASN
1	B	259	GLN
1	B	360	ASN
1	B	403	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 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	$\# Z > 2$
3	DHB	B	502	-	11,11,11	0.84	0	15,15,15	0.76	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
3	DHB	B	502	-	-	1/4/4/4	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	502	DHB	O2-C-C1-C2

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](#)

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	399/418 (95%)	0.44	12 (3%) 52 48	26, 76, 134, 161	0
1	B	404/418 (96%)	-0.08	5 (1%) 76 73	21, 44, 103, 150	0
All	All	803/836 (96%)	0.18	17 (2%) 63 59	21, 59, 126, 161	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	138	HIS	4.0
1	A	31	ALA	3.8
1	B	405	THR	3.6
1	A	18	PHE	3.6
1	A	262	PHE	3.4
1	A	200	GLY	2.9
1	A	231	LYS	2.7
1	A	215	ALA	2.7
1	A	33	ILE	2.6
1	A	2	ALA	2.6
1	A	40	VAL	2.2
1	A	139	GLY	2.2
1	B	363	TRP	2.2
1	B	399	VAL	2.1
1	B	137	GLU	2.1
1	A	255	LYS	2.1
1	A	206	TRP	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	DHB	B	502	11/11	0.95	0.08	26,27,30,32	0
2	FE2	A	501	1/1	0.96	0.19	32,32,32,32	0
2	FE2	B	501	1/1	0.99	0.01	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.