



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:17 PM GMT

PDB ID : 3WR3  
Title : Crystal structure of the anaerobic DesB-gallate complex  
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Deposited on : 2014-02-13  
Resolution : 2.50 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](mailto: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.

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

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

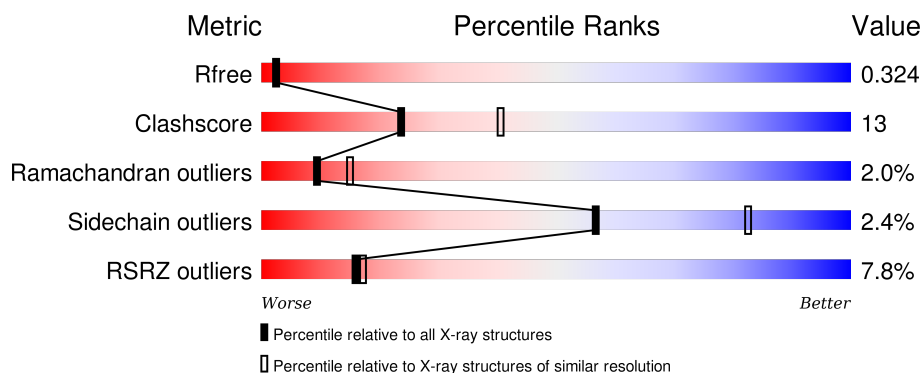
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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

Mol	Chain	Length	Quality of chain
1	A	418	<div> <div>10%</div> <div>62%</div> <div>32%</div> <div>• •</div> </div>
1	B	418	<div> <div>5%</div> <div>73%</div> <div>23%</div> <div>•</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6463 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.

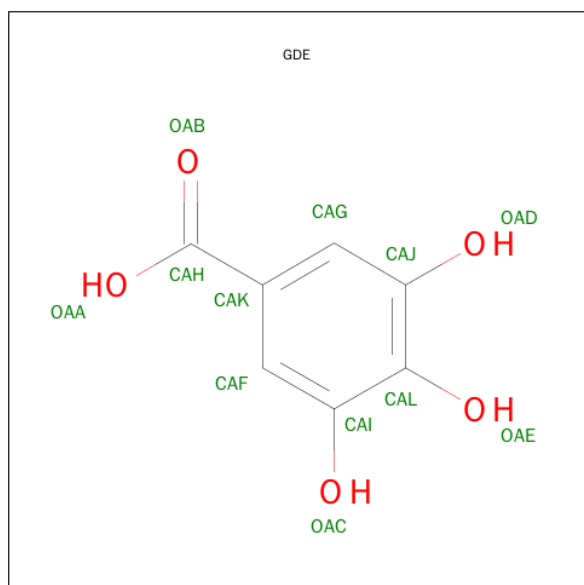
- Molecule 1 is a protein called Gallate dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	4	0	0
			3229	2074	554	587	14			
1	B	405	Total	C	N	O	S	0	0	0
			3220	2068	553	585	14			

- 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		

- Molecule 3 is 3,4,5-TRIHIDROXYBENZOIC ACID (three-letter code: GDE) (formula: C<sub>7</sub>H<sub>6</sub>O<sub>5</sub>).

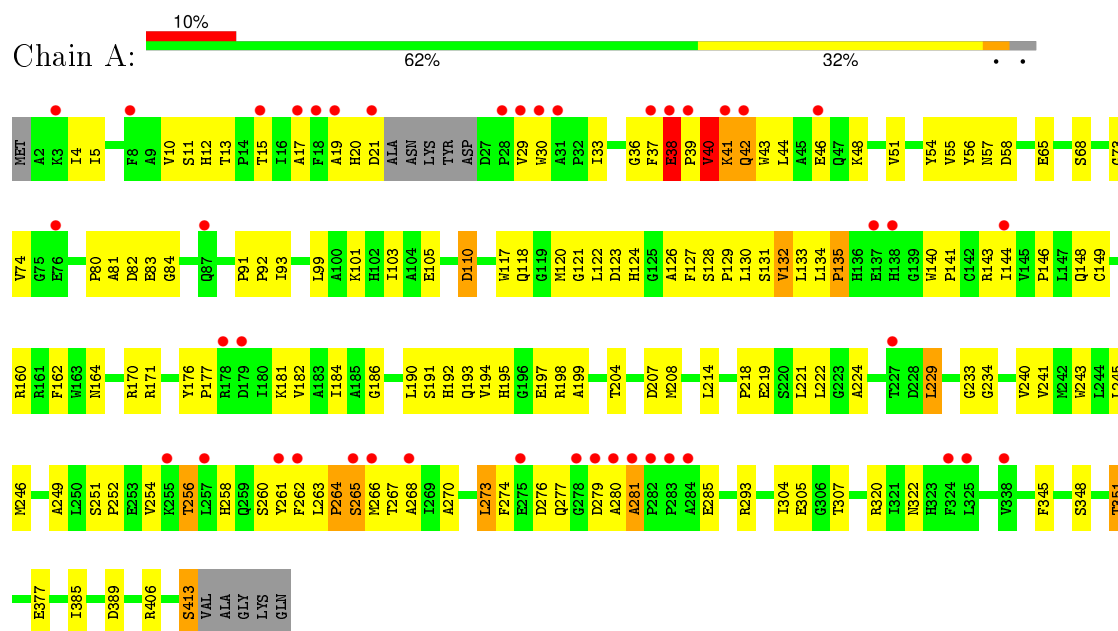


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	7	5		

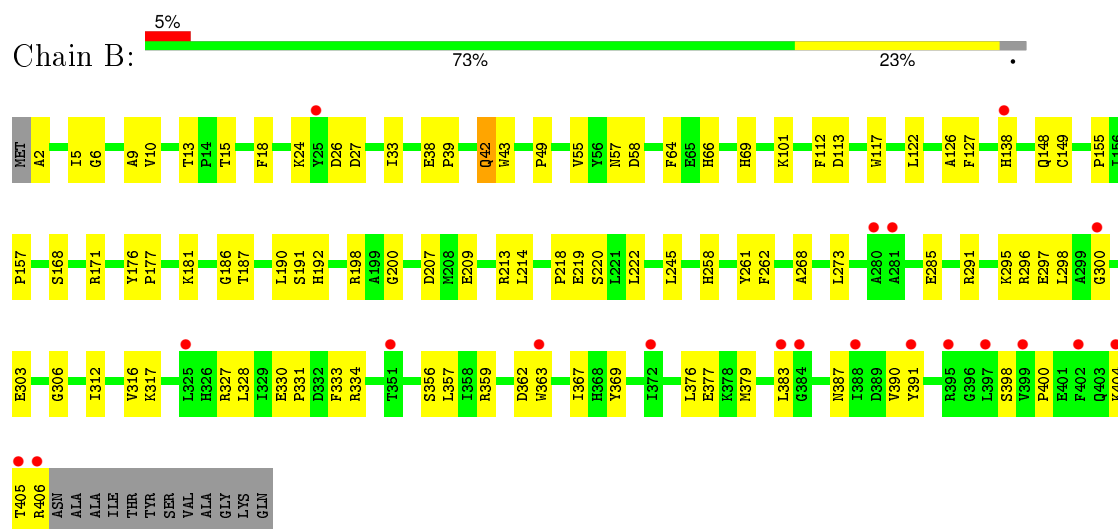
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: 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, $\alpha$ , $\beta$ , $\gamma$	58.35Å 64.02Å 118.48Å 90.00° 97.14° 90.00°	Depositor
Resolution (Å)	19.70 – 2.50 19.70 – 2.50	Depositor EDS
% Data completeness (in resolution range)	91.1 (19.70-2.50) 92.7 (19.70-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.36 (at 2.50Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1539)	Depositor
R, $R_{free}$	0.231 , 0.313 0.246 , 0.324	Depositor DCC
$R_{free}$ test set	1403 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	45.7	Xtriage
Anisotropy	0.988	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 40.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 28054 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6463	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.45	0/3323	0.64	2/4507 (0.0%)
1	B	0.50	0/3315	0.62	0/4496
All	All	0.47	0/6638	0.63	2/9003 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	263	LEU	CA-CB-CG	6.83	131.02	115.30
1	A	41	LYS	N-CA-C	-5.37	96.50	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	38	GLU	Peptide
1	A	40	VAL	Peptide

## 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	3229	0	3116	117	0
1	B	3220	0	3106	64	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	12	0	3	1	0
All	All	6463	0	6225	167	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 13.

All (167) 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:160:ARG:O	1:A:164:ASN:ND2	2.09	0.86
1:A:58:ASP:H	1:A:148:GLN:HE21	1.26	0.82
1:B:58:ASP:H	1:B:148:GLN:HE21	1.27	0.82
1:A:40:VAL:HA	1:A:43:TRP:H	1.45	0.82
1:A:38:GLU:HB3	1:A:41:LYS:HE2	1.61	0.81
1:A:20:HIS:CD2	1:A:132:VAL:HG23	2.16	0.81
1:A:170:ARG:NH2	1:A:279:ASP:OD1	2.15	0.80
1:A:171:ARG:NH2	1:A:285:GLU:O	2.19	0.76
1:A:40:VAL:HG23	1:A:43:TRP:HB3	1.69	0.74
1:B:377:GLU:HG3	1:B:387:ASN:HD21	1.56	0.71
1:A:20:HIS:HD2	1:A:132:VAL:HG23	1.57	0.70
1:A:82:ASP:OD2	1:B:334:ARG:NH2	2.26	0.69
1:A:13:THR:HG22	1:A:15:THR:H	1.58	0.69
1:A:39:PRO:HG2	1:A:258:HIS:ND1	2.09	0.68
1:A:170:ARG:HB2	1:A:249:ALA:HA	1.76	0.66
1:A:219:GLU:HA	1:A:222:LEU:HG	1.75	0.66
1:A:261:TYR:HE1	1:A:268:ALA:HB1	1.61	0.66
1:A:41:LYS:HD2	1:A:133:LEU:HA	1.78	0.65
1:A:265:SER:HA	1:B:406:ARG:HH22	1.61	0.65
1:A:99:LEU:O	1:A:103:ILE:HD12	1.98	0.64
1:A:54:TYR:HB2	1:A:144:ILE:HD11	1.79	0.63
1:A:40:VAL:HG22	1:A:44:LEU:HG	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:PRO:HG2	1:B:258:HIS:CE1	2.32	0.63
1:A:41:LYS:HD3	1:A:41:LYS:N	2.12	0.63
1:A:37:PHE:HB3	1:A:41:LYS:NZ	2.14	0.61
1:A:51:VAL:HG12	1:A:143:ARG:HB2	1.82	0.60
1:A:252:PRO:HG2	1:A:280:ALA:HB2	1.84	0.60
1:A:17:ALA:O	1:A:21:ASP:HB3	2.02	0.59
1:A:4:ILE:HD11	1:A:184:ILE:HG13	1.85	0.59
1:A:193:GLN:NE2	1:A:198:ARG:HB3	2.18	0.58
1:B:356:SER:HA	1:B:359:ARG:HG3	1.85	0.58
1:A:74:VAL:HG23	1:A:118:GLN:HG3	1.85	0.58
1:B:26:ASP:OD1	1:B:27:ASP:N	2.37	0.58
1:A:191:SER:HB3	1:A:207:ASP:OD1	2.04	0.57
1:B:356:SER:OG	1:B:359:ARG:NH1	2.38	0.57
1:B:191:SER:HB3	1:B:207:ASP:OD1	2.05	0.57
1:A:413:SER:HG	1:B:18:PHE:HE2	1.51	0.57
1:A:251:SER:OG	1:A:276:ASP:OD1	2.23	0.56
1:A:256:THR:HG22	1:A:258:HIS:O	2.05	0.56
1:A:73:GLY:HA3	1:A:146:PRO:HG2	1.85	0.56
1:A:29:VAL:HG23	1:A:30:TRP:CD1	2.40	0.56
1:B:331:PRO:HA	1:B:334:ARG:HB2	1.88	0.56
1:A:19:ALA:HB1	1:A:30:TRP:CH2	2.39	0.56
1:A:258:HIS:CE1	1:A:260:SER:HB3	2.41	0.56
1:A:51:VAL:HG23	1:A:182:VAL:HG13	1.88	0.56
1:A:101:LYS:NZ	1:A:105:GLU:OE2	2.32	0.56
1:A:80:PRO:HG3	1:A:127:PHE:HE2	1.71	0.56
1:A:184:ILE:HD13	1:A:249:ALA:HB3	1.86	0.56
1:A:176:TYR:CD1	1:A:177:PRO:HD2	2.40	0.56
1:B:69:HIS:CD2	1:B:157:PRO:HD3	2.41	0.55
1:A:38:GLU:CB	1:A:41:LYS:HE2	2.36	0.55
1:A:129:PRO:O	1:A:133:LEU:HG	2.06	0.55
1:A:214:LEU:O	1:A:218:PRO:HG3	2.05	0.55
1:A:91:PRO:HG2	1:A:140:TRP:NE1	2.23	0.54
1:B:261:TYR:CE1	1:B:268:ALA:HB1	2.43	0.54
1:A:42:GLN:O	1:A:46:GLU:HG3	2.07	0.53
1:A:124:HIS:HB2	1:A:128:SER:OG	2.08	0.53
1:A:110:ASP:OD2	1:A:293:ARG:NH1	2.42	0.53
1:B:112:PHE:HE2	1:B:168:SER:HG	1.56	0.52
1:A:224:ALA:HB1	1:A:229:LEU:HD13	1.91	0.52
1:A:37:PHE:HB3	1:A:41:LYS:HZ1	1.74	0.52
1:B:300:GLY:O	1:B:303:GLU:HG2	2.09	0.52
1:A:33:ILE:HG22	1:A:262:PHE:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:TRP:O	1:A:33:ILE:HG12	2.10	0.51
1:A:186:GLY:H	1:A:246:MET:HG3	1.75	0.51
1:A:190:LEU:HD11	1:A:243:TRP:CH2	2.45	0.51
1:A:122:LEU:HD12	1:A:127:PHE:HE1	1.75	0.51
1:B:295:LYS:HZ3	1:B:298:LEU:HB2	1.76	0.51
1:A:5:ILE:HG23	1:A:181:LYS:HB3	1.93	0.50
1:A:93:ILE:HG13	1:A:140:TRP:HB2	1.92	0.50
1:A:320:ARG:HH21	1:A:351:THR:HG23	1.74	0.50
1:B:58:ASP:N	1:B:148:GLN:HE21	2.05	0.50
1:A:345:PHE:O	1:A:348:SER:OG	2.22	0.50
1:A:148:GLN:NE2	1:A:149:CYS:O	2.44	0.49
1:B:58:ASP:H	1:B:148:GLN:NE2	2.02	0.49
1:A:170:ARG:HB2	1:A:249:ALA:CA	2.40	0.49
1:B:363:TRP:CE2	1:B:390:VAL:HG22	2.48	0.49
1:A:197:GLU:HB2	1:B:306:GLY:HA3	1.95	0.49
1:A:134:LEU:HD22	1:A:141:PRO:HG2	1.95	0.48
1:A:83:GLU:OE1	1:A:123:ASP:HB2	2.13	0.48
1:A:385:ILE:O	1:A:385:ILE:HG13	2.12	0.48
1:B:377:GLU:HG3	1:B:387:ASN:ND2	2.27	0.48
1:B:171:ARG:NH2	1:B:285:GLU:O	2.39	0.48
1:B:331:PRO:HA	1:B:334:ARG:HE	1.79	0.47
1:B:43:TRP:CD2	1:B:273:LEU:HD11	2.49	0.47
1:A:254:VAL:HG21	1:A:274:PHE:CD1	2.49	0.47
1:A:377:GLU:OE2	3:A:502:GDE:OAC	2.33	0.47
1:B:55:VAL:HG21	1:B:245:LEU:HD22	1.95	0.47
1:B:214:LEU:O	1:B:218:PRO:HG3	2.15	0.47
1:A:40:VAL:HG21	1:A:273:LEU:HD21	1.97	0.46
1:A:20:HIS:HD1	1:A:20:HIS:C	2.17	0.46
1:B:33:ILE:HG13	1:B:262:PHE:CG	2.51	0.46
1:A:190:LEU:HD13	1:A:270:ALA:HB3	1.96	0.46
1:B:376:LEU:O	1:B:379:MET:N	2.49	0.46
1:A:265:SER:HA	1:B:406:ARG:NH2	2.29	0.46
1:B:387:ASN:O	1:B:391:TYR:HD2	1.99	0.45
1:B:38:GLU:O	1:B:42:GLN:HG3	2.16	0.45
1:A:130:LEU:C	1:A:132:VAL:H	2.18	0.45
1:A:65:GLU:O	1:B:117:TRP:HA	2.16	0.45
1:B:328:LEU:HD23	1:B:333:PHE:HD2	1.81	0.45
1:B:13:THR:HG22	1:B:15:THR:H	1.81	0.45
1:A:261:TYR:CE1	1:A:268:ALA:HB1	2.47	0.45
1:B:9:ALA:HA	1:B:186:GLY:O	2.16	0.45
1:A:38:GLU:CB	1:A:41:LYS:HG2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:LEU:CD1	1:A:270:ALA:HB3	2.47	0.45
1:A:55:VAL:HB	1:A:186:GLY:HA2	1.98	0.45
1:B:357:LEU:HD23	1:B:362:ASP:HB3	1.99	0.45
1:A:204:THR:O	1:A:208:MET:HG2	2.16	0.45
1:A:322:ASN:HB3	1:B:64:PHE:CZ	2.51	0.45
1:B:219:GLU:HA	1:B:222:LEU:HG	1.98	0.45
1:A:305:GLU:O	1:B:198:ARG:HD2	2.17	0.44
1:A:91:PRO:HA	1:A:92:PRO:HD2	1.82	0.44
1:A:117:TRP:HB2	1:A:120:MET:HE2	2.00	0.44
1:A:162:PHE:O	1:A:245:LEU:HD12	2.18	0.44
1:B:24:LYS:C	1:B:26:ASP:H	2.21	0.44
1:A:105:GLU:HG2	1:B:155:PRO:HG3	2.00	0.44
1:A:264:PRO:HB2	1:A:265:SER:H	1.65	0.44
1:A:84:GLY:C	1:B:334:ARG:HH12	2.22	0.43
1:A:10:VAL:HG22	1:A:11:SER:O	2.18	0.43
1:A:134:LEU:HA	1:A:135:PRO:HD2	1.80	0.43
1:A:40:VAL:O	1:A:40:VAL:HG22	2.19	0.43
1:A:54:TYR:OH	1:A:126:ALA:HA	2.18	0.43
1:B:177:PRO:HD3	1:B:291:ARG:HD3	1.99	0.43
1:A:4:ILE:HG23	1:A:276:ASP:OD2	2.18	0.43
1:A:101:LYS:O	1:A:105:GLU:HG3	2.19	0.43
1:B:177:PRO:HB3	1:B:291:ARG:HH11	1.83	0.43
1:A:68:SER:OG	1:B:66:HIS:NE2	2.48	0.43
1:A:40:VAL:CG2	1:A:273:LEU:HD21	2.48	0.43
1:A:54:TYR:CE2	1:A:56:TYR:HB3	2.53	0.43
1:A:304:ILE:HG22	1:A:307:THR:OG1	2.19	0.43
1:A:81:ALA:HB2	1:A:121:GLY:C	2.39	0.43
1:B:5:ILE:HG13	1:B:6:GLY:N	2.33	0.43
1:B:101:LYS:HB3	1:B:101:LYS:HE3	1.77	0.43
1:B:331:PRO:HA	1:B:334:ARG:NE	2.34	0.43
1:A:192:HIS:N	1:A:233:GLY:O	2.52	0.43
1:B:317:LYS:HE3	1:B:369:TYR:O	2.19	0.43
1:A:20:HIS:ND1	1:A:20:HIS:C	2.72	0.43
1:A:38:GLU:HB3	1:A:41:LYS:HG2	2.00	0.42
1:B:312:ILE:O	1:B:316:VAL:HG23	2.19	0.42
1:A:4:ILE:HD12	1:A:182:VAL:O	2.19	0.42
1:A:117:TRP:HH2	1:A:122:LEU:HD21	1.84	0.42
1:A:218:PRO:O	1:A:221:LEU:HB2	2.19	0.42
1:A:170:ARG:HG3	1:A:249:ALA:O	2.19	0.42
1:A:171:ARG:NH1	1:A:285:GLU:O	2.53	0.42
1:A:84:GLY:O	1:B:334:ARG:NH1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:HIS:HB3	1:A:129:PRO:HG3	2.00	0.42
1:B:191:SER:O	1:B:192:HIS:HB3	2.19	0.42
1:B:2:ALA:HB2	1:B:176:TYR:O	2.20	0.42
1:B:327:ARG:NE	1:B:330:GLU:OE2	2.37	0.42
1:A:267:THR:HG22	1:A:268:ALA:N	2.35	0.41
1:B:398:SER:OG	1:B:400:PRO:HD2	2.20	0.41
1:A:279:ASP:HB3	1:A:281:ALA:O	2.20	0.41
1:B:69:HIS:O	1:B:149:CYS:HA	2.20	0.41
1:A:54:TYR:CE2	1:A:126:ALA:HA	2.55	0.41
1:B:122:LEU:HB3	1:B:126:ALA:HB3	2.02	0.41
1:A:195:HIS:HB2	1:B:367:ILE:HD12	2.03	0.41
1:A:130:LEU:O	1:A:132:VAL:N	2.54	0.41
1:A:33:ILE:CG2	1:A:262:PHE:HB3	2.51	0.41
1:B:49:PRO:HA	1:B:181:LYS:HB2	2.03	0.41
1:A:194:VAL:HG11	1:A:266:MET:CE	2.51	0.41
1:A:36:GLY:O	1:A:258:HIS:CE1	2.74	0.41
1:A:406:ARG:NH2	1:B:200:GLY:O	2.51	0.40
1:B:209:GLU:OE2	1:B:213:ARG:NH1	2.53	0.40
1:B:122:LEU:HD12	1:B:127:PHE:CE2	2.56	0.40
1:B:10:VAL:O	1:B:187:THR:HA	2.21	0.40
1:B:297:GLU:O	1:B:298:LEU:HD23	2.21	0.40
1:A:192:HIS:O	1:A:234:GLY:HA3	2.22	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	403/418 (96%)	361 (90%)	30 (7%)	12 (3%)	5	7
1	B	403/418 (96%)	365 (91%)	34 (8%)	4 (1%)	19	34
All	All	806/836 (96%)	726 (90%)	64 (8%)	16 (2%)	9	15

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	GLU
1	A	265	SER
1	A	281	ALA
1	B	296	ARG
1	A	40	VAL
1	A	131	SER
1	A	42	GLN
1	A	264	PRO
1	A	199	ALA
1	A	256	THR
1	B	190	LEU
1	B	404	LYS
1	A	351	THR
1	B	138	HIS
1	A	135	PRO
1	A	240	VAL

### 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	332/340 (98%)	322 (97%)	10 (3%)	48	76
1	B	331/340 (97%)	325 (98%)	6 (2%)	66	88
All	All	663/680 (98%)	647 (98%)	16 (2%)	57	82

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LYS
1	A	57	ASN
1	A	110	ASP
1	A	132	VAL
1	A	229	LEU
1	A	241	VAL
1	A	273	LEU

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Mol	Chain	Res	Type
1	A	277	GLN
1	A	389	ASP
1	A	413	SER
1	B	42	GLN
1	B	57	ASN
1	B	113	ASP
1	B	220	SER
1	B	383	LEU
1	B	405	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	A	277	GLN
1	B	148	GLN
1	B	387	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 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	GDE	A	502	-	9,12,12	0.87	1 (11%)	12,17,17	1.20	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	GDE	A	502	-	-	0/0/4/4	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	GDE	CAJ-CAL	-2.23	1.37	1.39

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	GDE	1	0

## 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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	407/418 (97%)	0.58	43 (10%) 8 8	37, 65, 94, 119	4 (0%)
1	B	405/418 (96%)	0.28	20 (4%) 33 38	34, 50, 82, 101	1 (0%)
All	All	812/836 (97%)	0.43	63 (7%) 16 17	34, 57, 89, 119	5 (0%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	280	ALA	7.5
1	A	30	TRP	7.3
1	A	18	PHE	6.0
1	A	281	ALA	5.4
1	A	17	ALA	4.4
1	A	46	GLU	4.4
1	A	42	GLN	4.4
1	B	406	ARG	4.3
1	A	19	ALA	4.2
1	A	262	PHE	4.2
1	A	28	PRO	4.0
1	B	388	ILE	3.7
1	A	41	LYS	3.7
1	A	282	PRO	3.6
1	A	261	TYR	3.6
1	B	391	TYR	3.5
1	A	278	GLY	3.4
1	A	266	MET	3.4
1	B	384	GLY	3.3
1	B	383	LEU	3.3
1	B	397	LEU	3.1
1	A	179	ASP	3.1
1	B	405	THR	3.1
1	A	39	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	138	HIS	3.0
1	A	138	HIS	2.9
1	A	38	GLU	2.9
1	A	87	GLN	2.9
1	A	29	VAL	2.9
1	B	395	ARG	2.9
1	B	281	ALA	2.8
1	A	3	LYS	2.7
1	A	257	LEU	2.7
1	B	25	TYR	2.7
1	A	76	GLU	2.7
1	B	280	ALA	2.6
1	A	279	ASP	2.6
1	A	15	THR	2.5
1	A	283	PRO	2.5
1	A	255	LYS	2.4
1	A	325	LEU	2.4
1	A	265	SER	2.4
1	A	324	PHE	2.4
1	B	300	GLY	2.3
1	B	363	TRP	2.3
1	A	137	GLU	2.3
1	A	338	VAL	2.3
1	A	8	PHE	2.3
1	A	178	ARG	2.2
1	A	31	ALA	2.2
1	A	37	PHE	2.2
1	B	404	LYS	2.2
1	A	275	GLU	2.2
1	A	268	ALA	2.2
1	A	21	ASP	2.1
1	A	227	THR	2.1
1	B	351	THR	2.1
1	B	399	VAL	2.1
1	B	325	LEU	2.1
1	A	284	ALA	2.1
1	B	372	ILE	2.0
1	A	144	ILE	2.0
1	B	402	PHE	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GDE	A	502	12/12	0.91	0.16	-0.04	38,41,44,46	0
2	FE2	B	501	1/1	0.98	0.10	-	47,47,47,47	0
2	FE2	A	501	1/1	0.96	0.04	-	72,72,72,72	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.