



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:25 AM GMT

PDB ID : 2QMA
Title : Crystal structure of glutamate decarboxylase domain of diaminobutyrate-pyruvate transaminase and L-2,4-diaminobutyrate decarboxylase from *Vibrio parahaemolyticus*
Authors : Osipiuk, J.; Sather, A.; Gu, M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2007-07-14
Resolution : 1.81 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) : 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

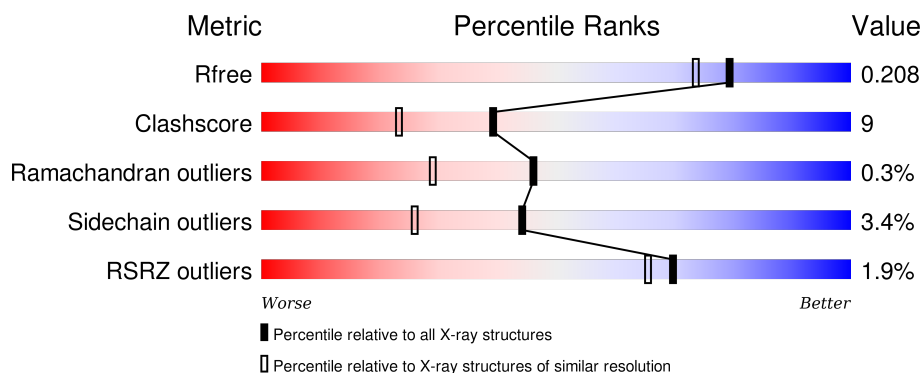
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

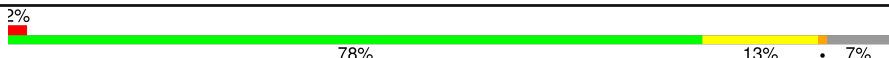

The reported resolution of this entry is 1.81 Å.

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	5422 (1.84-1.80)
Clashscore	102246	6347 (1.84-1.80)
Ramachandran outliers	100387	6276 (1.84-1.80)
Sidechain outliers	100360	6276 (1.84-1.80)
RSRZ outliers	91569	5439 (1.84-1.80)

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.

Mol	Chain	Length	Quality of chain
1	A	497	
1	B	497	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	B	962	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7843 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 Diaminobutyrate-pyruvate transaminase and L-2,4-diaminobutyrate decarboxylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	463	Total	C	N	O	S	Se	0	7	0
			3580	2269	595	690	5	21			
1	B	459	Total	C	N	O	S	Se	0	14	0
			3582	2273	591	690	5	23			

There are 44 discrepancies between the modelled and reference sequences:

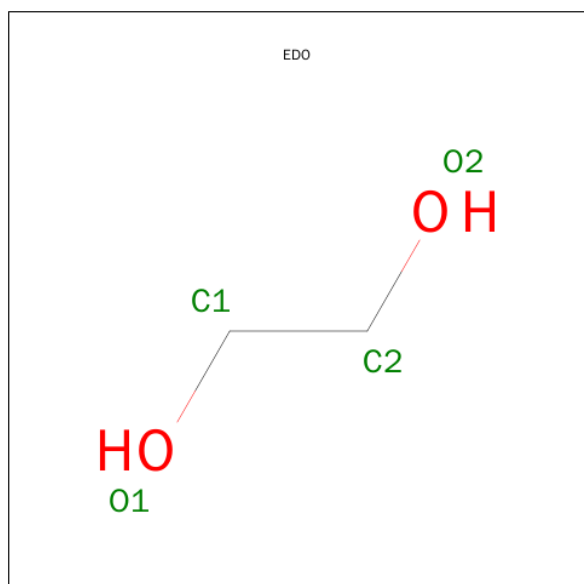
Chain	Residue	Modelled	Actual	Comment	Reference
A	462	SER	-	CLONING ARTIFACT	UNP Q87NC6
A	463	ASN	-	CLONING ARTIFACT	UNP Q87NC6
A	464	ALA	-	CLONING ARTIFACT	UNP Q87NC6
A	499	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	506	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	520	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	574	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	582	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	590	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	634	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	637	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	683	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	691	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	701	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	744	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	750	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	752	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	843	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	845	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	856	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	870	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
A	952	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	462	SER	-	CLONING ARTIFACT	UNP Q87NC6
B	463	ASN	-	CLONING ARTIFACT	UNP Q87NC6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	464	ALA	-	CLONING ARTIFACT	UNP Q87NC6
B	499	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	506	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	520	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	574	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	582	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	590	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	634	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	637	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	683	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	691	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	701	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	744	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	750	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	752	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	843	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	845	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	856	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	870	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6
B	952	MSE	MET	MODIFIED RESIDUE	UNP Q87NC6

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	343	Total	O	0	2
			343	343		
3	B	328	Total	O	0	3
			330	330		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.43Å 122.78Å 64.29Å 90.00° 95.74° 90.00°	Depositor
Resolution (Å)	37.67 – 1.81 37.68 – 1.81	Depositor EDS
% Data completeness (in resolution range)	95.8 (37.67-1.81) 95.4 (37.68-1.81)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.181 , 0.209 0.181 , 0.208	Depositor DCC
R_{free} test set	4331 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.140	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.1	EDS
Estimated twinning fraction	0.267 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 86394 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7843	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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, 0.375 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: EDO

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	0.77	0/3648	0.76	4/4906 (0.1%)
1	B	0.76	0/3667	0.78	7/4926 (0.1%)
All	All	0.76	0/7315	0.77	11/9832 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	783[A]	ASP	CB-CA-C	6.56	123.53	110.40
1	B	783[B]	ASP	CB-CA-C	6.56	123.53	110.40
1	B	783[A]	ASP	CB-CG-OD2	-6.06	112.85	118.30
1	B	783[B]	ASP	CB-CG-OD2	-6.06	112.85	118.30
1	B	787	LEU	CA-CB-CG	5.97	129.04	115.30
1	A	809	LEU	CA-CB-CG	5.78	128.58	115.30
1	A	623	ILE	CG1-CB-CG2	5.57	123.65	111.40
1	A	623	ILE	CB-CA-C	-5.46	100.68	111.60
1	A	691	MSE	CG-SE-CE	-5.32	87.19	98.90
1	B	783[A]	ASP	CB-CG-OD1	5.19	122.97	118.30
1	B	783[B]	ASP	CB-CG-OD1	5.19	122.97	118.30

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	3580	0	3575	77	0
1	B	3582	0	3601	64	0
2	A	4	0	6	0	0
2	B	4	0	6	0	0
3	A	343	0	0	19	0
3	B	330	0	0	12	0
All	All	7843	0	7188	128	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 9.

All (128) 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:844:THR:HG22	1:A:845:MSE:CE	1.52	1.40
1:B:633[B]:GLN:HG2	3:B:1219[B]:HOH:O	1.34	1.23
1:A:548:VAL:HG13	3:A:1235:HOH:O	1.48	1.13
3:A:1284:HOH:O	1:B:534:LEU:HG	1.56	1.05
1:A:844:THR:HG22	1:A:845:MSE:HE2	1.05	1.02
3:A:1296:HOH:O	1:B:548:VAL:HG13	1.59	1.02
1:A:844:THR:CG2	1:A:845:MSE:CE	2.38	1.00
1:A:683:MSE:HE2	3:A:1149:HOH:O	1.61	1.00
1:A:605:LYS:HE3	3:A:1284:HOH:O	1.61	0.99
1:A:845:MSE:HE1	1:A:853:LEU:HD11	1.47	0.96
1:B:785:HIS:HD2	1:B:794:CYS:H	1.16	0.93
1:B:731:ASP:OD2	1:B:921:THR:HG21	1.70	0.92
1:B:674:PHE:HD1	1:B:678:LYS:HD2	1.34	0.92
1:A:634[B]:MSE:HE3	1:A:808:LEU:HD11	1.59	0.83
1:A:617:SER:HB3	3:A:1226:HOH:O	1.81	0.81
1:B:582[B]:MSE:HE3	1:B:586:LEU:HD11	1.64	0.80
1:A:844:THR:CG2	1:A:845:MSE:HE2	2.01	0.79
1:B:633[B]:GLN:CG	3:B:1219[B]:HOH:O	2.03	0.78
1:A:785:HIS:HD2	1:A:794:CYS:H	1.32	0.78
1:B:674:PHE:CD1	1:B:678:LYS:HD2	2.19	0.77
1:B:944[A]:GLU:HG2	3:B:1283:HOH:O	1.84	0.76
1:B:894:HIS:HD2	1:B:897:ALA:H	1.34	0.76
1:B:633[B]:GLN:CB	3:B:1219[B]:HOH:O	2.33	0.75
1:A:691:MSE:HE1	1:A:713:ALA:HB2	1.68	0.75
1:B:785:HIS:CD2	1:B:794:CYS:H	2.05	0.74
1:A:634[B]:MSE:CE	1:A:805:PHE:HE1	2.00	0.73
1:B:479:GLU:O	1:B:482:LYS:HE2	1.88	0.73
1:A:845:MSE:HE1	1:A:853:LEU:CD1	2.17	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:785:HIS:HD2	1:B:794:CYS:N	1.86	0.71
1:B:650:HIS:HE1	1:B:656:GLY:O	1.76	0.69
1:B:617:SER:HB3	3:B:1165:HOH:O	1.92	0.68
1:A:844:THR:HG22	1:A:845:MSE:HE3	1.65	0.68
1:B:544:VAL:O	1:B:548:VAL:HG22	1.94	0.67
3:A:1043:HOH:O	1:B:483:HIS:HD2	1.78	0.67
1:A:907:ARG:HD2	1:A:920:GLU:HB3	1.76	0.66
1:A:582[B]:MSE:HE1	1:B:548:VAL:CG2	2.25	0.65
1:B:582[B]:MSE:HE3	1:B:586:LEU:CD1	2.26	0.65
1:A:634[B]:MSE:HE1	1:A:805:PHE:CE1	2.32	0.65
1:A:548:VAL:HG23	1:B:582[A]:MSE:HE1	1.79	0.65
1:A:634[B]:MSE:CE	1:A:805:PHE:CE1	2.82	0.62
1:A:552:VAL:CG1	1:B:582[B]:MSE:HE1	2.29	0.62
1:A:554:LYS:HD2	3:A:1276:HOH:O	1.98	0.62
1:B:497:SER:OG	1:B:542:LYS:NZ	2.34	0.61
1:A:682:TRP:CD1	1:B:637:MSE:HE1	2.35	0.61
1:B:742:ALA:HB2	1:B:752:MSE:HE1	1.82	0.61
1:A:623:ILE:HD11	1:A:634[B]:MSE:HE1	1.81	0.60
1:B:590:MSE:HB2	1:B:834[B]:LYS:HB3	1.83	0.60
1:A:476:PRO:HB2	1:A:479:GLU:HG2	1.84	0.59
1:A:654:LYS:HD2	3:A:1136:HOH:O	2.01	0.59
1:B:894:HIS:CD2	1:B:897:ALA:H	2.19	0.59
1:A:552:VAL:HG12	1:B:582[B]:MSE:HE1	1.84	0.59
1:B:731:ASP:CG	1:B:921:THR:HG21	2.22	0.58
1:A:548:VAL:CG2	1:B:582[A]:MSE:HE1	2.34	0.58
1:A:605:LYS:CE	3:A:1284:HOH:O	2.33	0.57
1:A:623:ILE:CD1	1:A:634[A]:MSE:HE1	2.35	0.57
1:A:634[B]:MSE:HE2	1:A:797:LEU:HD23	1.87	0.56
1:B:659:ASP:HB2	3:B:1107:HOH:O	2.04	0.56
1:B:633[B]:GLN:OE1	1:B:679:SER:HA	2.05	0.56
1:A:563:ASP:OD2	1:A:949:LYS:HE3	2.06	0.55
1:B:673:HIS:CD2	1:B:675:THR:HG23	2.42	0.53
1:A:582[B]:MSE:HG3	1:B:506:MSE:SE	2.58	0.53
1:B:509:VAL:O	1:B:513:VAL:HG23	2.08	0.52
1:B:623:ILE:HG21	1:B:634[B]:MSE:HE1	1.92	0.51
1:B:923:VAL:O	1:B:926:LYS:HG2	2.10	0.51
1:A:582[B]:MSE:HE1	1:B:548:VAL:HG21	1.92	0.51
1:A:631:SER:HB3	1:A:781:SER:HB3	1.93	0.51
1:A:844:THR:CG2	1:A:845:MSE:HE3	2.30	0.50
1:A:683:MSE:CE	3:A:1149:HOH:O	2.36	0.50
1:A:835:ARG:HB3	1:B:792:ILE:HG22	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:623:ILE:HD13	1:A:634[A]:MSE:HE1	1.95	0.48
1:B:924:ASP:HB2	3:B:1145:HOH:O	2.13	0.48
1:A:623:ILE:HD11	1:A:634[B]:MSE:CE	2.43	0.48
1:A:604:GLN:HE21	1:A:608:ASN:ND2	2.12	0.48
1:A:891:ARG:NH1	1:A:926:LYS:HD2	2.29	0.48
1:A:623:ILE:HD11	1:A:634[A]:MSE:HE1	1.96	0.48
1:B:628:GLY:N	1:B:783[B]:ASP:OD1	2.46	0.48
1:A:515:ALA:HB1	1:A:914:GLY:O	2.13	0.48
1:A:490:LEU:N	1:A:490:LEU:HD12	2.29	0.48
1:A:875:ASP:OD1	1:A:875:ASP:N	2.47	0.47
1:A:785:HIS:CD2	1:A:794:CYS:H	2.22	0.47
1:B:633[B]:GLN:OE1	1:B:682:TRP:HE3	1.98	0.47
1:A:785:HIS:HD2	1:A:794:CYS:N	2.05	0.46
1:A:565:ILE:HG23	1:A:565:ILE:O	2.15	0.46
1:B:574:MSE:N	1:B:575:PRO:HD2	2.31	0.46
1:A:574:MSE:N	1:A:575:PRO:HD2	2.31	0.46
1:B:650:HIS:HD2	3:B:1226:HOH:O	1.97	0.46
1:A:619:LYS:NZ	3:A:1293:HOH:O	2.49	0.46
1:A:634[B]:MSE:HE2	1:A:805:PHE:HE1	1.80	0.45
1:B:633[A]:GLN:HG3	1:B:634[A]:MSE:HE3	1.98	0.45
1:A:948:SER:O	1:A:952:MSE:HG2	2.17	0.45
1:A:833:THR:OG1	3:A:1280:HOH:O	2.21	0.45
1:A:637:MSE:HE1	3:A:1259:HOH:O	2.16	0.44
1:A:610:LEU:HD11	1:A:784:PHE:HZ	1.82	0.44
1:A:605:LYS:NZ	3:A:1284:HOH:O	2.46	0.44
1:B:482:LYS:HB3	3:B:1178:HOH:O	2.18	0.44
1:A:945:SER:O	1:A:949:LYS:HG3	2.17	0.44
1:A:604:GLN:HE21	1:A:608:ASN:HD21	1.64	0.44
1:A:674[B]:PHE:HD2	1:A:678:LYS:HD2	1.83	0.43
1:A:682:TRP:HD1	1:B:637:MSE:HE1	1.82	0.43
1:B:909:GLU:HG2	3:B:1262:HOH:O	2.18	0.43
3:A:1236:HOH:O	1:B:483:HIS:HE1	2.01	0.43
1:B:562:PRO:HB3	1:B:938:LEU:HD22	2.01	0.43
1:B:757:ALA:HA	1:B:783[B]:ASP:HB2	2.01	0.43
1:A:832:THR:O	1:A:833:THR:C	2.55	0.43
1:A:843:MSE:SE	1:B:548:VAL:HG21	2.69	0.43
1:B:621:ASP:HB2	1:B:802:LYS:HB3	2.01	0.42
1:A:727:ALA:HA	1:A:735:ILE:HG23	2.01	0.42
1:A:637:MSE:CE	3:A:1259:HOH:O	2.67	0.42
1:B:636:LEU:CB	1:B:683:MSE:HE1	2.49	0.42
1:B:650:HIS:CE1	1:B:656:GLY:O	2.65	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:573:LEU:HG	1:A:789:TYR:HB2	2.00	0.42
1:A:590:MSE:HG2	1:A:834:LYS:O	2.18	0.42
1:A:761:ALA:HB3	1:A:787:LEU:HD23	2.02	0.42
1:A:633:GLN:HG3	3:A:1149:HOH:O	2.19	0.42
1:A:573:LEU:HD22	1:B:483:HIS:HB3	2.02	0.42
1:A:631:SER:HB3	1:A:781:SER:CB	2.50	0.42
1:B:520:MSE:HE2	1:B:525:LEU:N	2.35	0.41
1:A:844:THR:C	1:A:845:MSE:HE2	2.39	0.41
1:A:552:VAL:HG12	1:B:582[B]:MSE:CE	2.49	0.41
1:A:623:ILE:HD11	1:A:634[B]:MSE:SE	2.70	0.41
1:B:752:MSE:HE3	1:B:777:ALA:HA	2.02	0.41
1:B:507:LYS:O	1:B:511:GLU:HG3	2.20	0.41
1:A:906:LEU:C	1:A:906:LEU:HD12	2.41	0.41
1:B:859:HIS:HD2	3:B:1288:HOH:O	2.03	0.41
1:A:613:LYS:HA	1:A:613:LYS:HD3	1.94	0.40
1:A:866:GLU:HB3	3:A:1233:HOH:O	2.21	0.40
1:B:650:HIS:CD2	3:B:1226:HOH:O	2.74	0.40
1:B:515:ALA:HB1	1:B:914:GLY:O	2.21	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	466/497 (94%)	449 (96%)	15 (3%)	2 (0%)	39	23
1	B	468/497 (94%)	454 (97%)	13 (3%)	1 (0%)	52	35
All	All	934/994 (94%)	903 (97%)	28 (3%)	3 (0%)	46	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	833	THR
1	B	596	ALA
1	A	565	ILE

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	390/391 (100%)	377 (97%)	13 (3%)	45	27
1	B	394/391 (101%)	379 (96%)	15 (4%)	40	21
All	All	784/782 (100%)	756 (96%)	28 (4%)	44	24

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	492	SER
1	A	612	ASP
1	A	623	ILE
1	A	688	LYS
1	A	691	MSE
1	A	783	ASP
1	A	800	ASN
1	A	810	HIS
1	A	875	ASP
1	A	885	LEU
1	A	895	GLU
1	A	906	LEU
1	A	907	ARG
1	B	490[A]	LEU
1	B	490[B]	LEU
1	B	548	VAL
1	B	612	ASP
1	B	739	ASP
1	B	781	SER
1	B	783[A]	ASP
1	B	783[B]	ASP

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Mol	Chain	Res	Type
1	B	800	ASN
1	B	906	LEU
1	B	907	ARG
1	B	909	GLU
1	B	913	ARG
1	B	921	THR
1	B	940	THR

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

Mol	Chain	Res	Type
1	A	530	ASN
1	A	608	ASN
1	A	732	HIS
1	A	785	HIS
1	A	859	HIS
1	B	483	HIS
1	B	650	HIS
1	B	673	HIS
1	B	696	ASN
1	B	785	HIS
1	B	859	HIS
1	B	894	HIS

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](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	EDO	A	961	-	3,3,3	0.58	0	2,2,2	0.17	0
2	EDO	B	962	-	3,3,3	0.73	0	2,2,2	0.47	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
2	EDO	A	961	-	-	0/1/1/1	0/0/0/0
2	EDO	B	962	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	444/497 (89%)	-0.33	9 (2%) 68 64	7, 18, 30, 39	0
1	B	440/497 (88%)	-0.37	8 (1%) 71 67	8, 18, 27, 33	0
All	All	884/994 (88%)	-0.35	17 (1%) 70 65	7, 18, 29, 39	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	810	HIS	4.6
1	B	475	ALA	4.5
1	B	674	PHE	4.1
1	A	833	THR	3.9
1	A	674[A]	PHE	3.8
1	B	490[A]	LEU	3.0
1	A	682	TRP	2.9
1	B	548	VAL	2.8
1	B	809	LEU	2.7
1	B	673	HIS	2.6
1	A	673	HIS	2.4
1	B	476	PRO	2.3
1	A	925	GLY	2.3
1	B	672	SER	2.2
1	A	490	LEU	2.2
1	A	832	THR	2.1
1	A	612	ASP	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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	EDO	B	962	4/4	0.93	0.09	2.60	25,27,28,29	0
2	EDO	A	961	4/4	0.96	0.11	-	24,24,24,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.