Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 10:36 PM EDT

PDB ID	:	3D3S
Title	:	Crystal structure of L-2,4-diaminobutyric acid acetyltransferase from Bordetella parapertussis
Authors	:	Kim, Y.; Volkart, L.; Jedrzejczak, R.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2008-05-12
Resolution	:	1.87 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 (i) 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 (i)) were used in the production of this report:

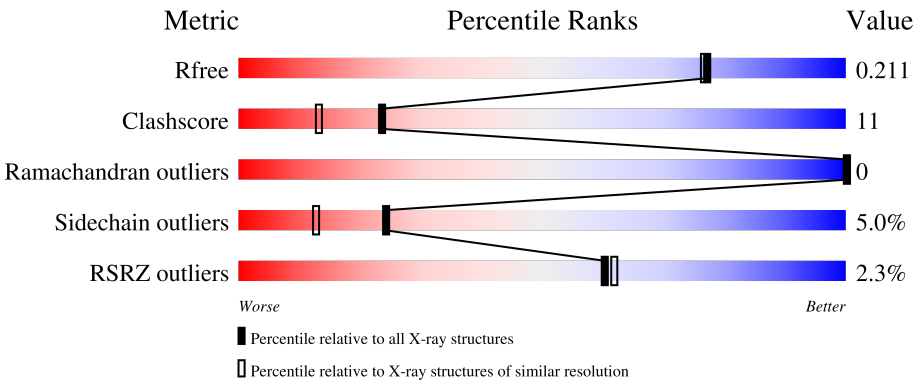
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 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}	130704	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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 <=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	189	3%, 66%, 17%, •, 16%
1	B	189	%, 64%, 13%, •, 19%
1	C	189	2%, 63%, 12%, •, 21%
1	D	189	2%, 65%, 11%, •, 23%

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	SO4	D	166	-	-	X	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5650 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.

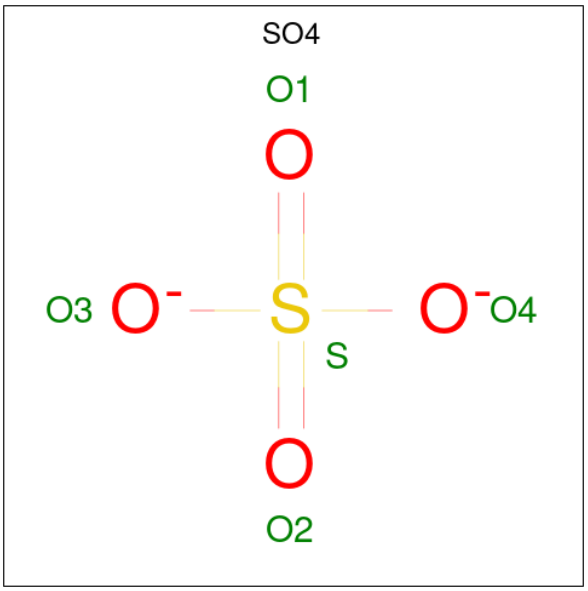
- Molecule 1 is a protein called L-2,4-diaminobutyric acid acetyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	159	1349	836	269	238	4	2	0	10	0
1	B	153	1333	830	269	227	4	3	0	12	0
1	C	149	1259	788	249	216	4	2	0	8	0
1	D	145	1232	767	250	209	4	2	0	8	0

There are 12 discrepancies between the modelled and reference sequences:

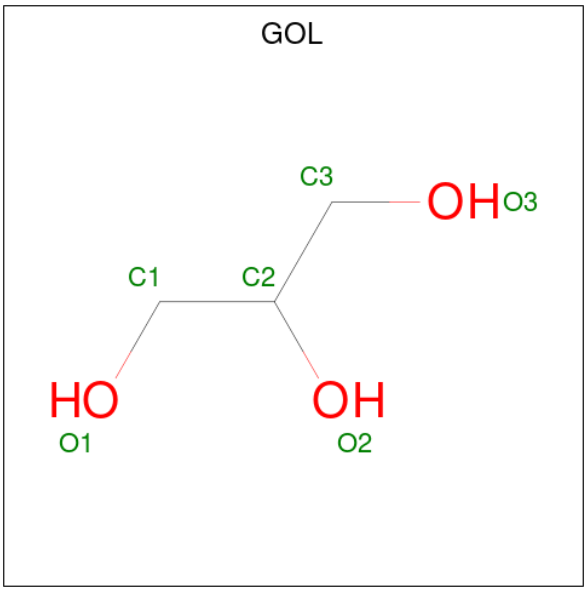
Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	SER	-	EXPRESSION TAG	UNP Q7W980
A	-23	ASN	-	EXPRESSION TAG	UNP Q7W980
A	-22	ALA	-	EXPRESSION TAG	UNP Q7W980
B	-24	SER	-	EXPRESSION TAG	UNP Q7W980
B	-23	ASN	-	EXPRESSION TAG	UNP Q7W980
B	-22	ALA	-	EXPRESSION TAG	UNP Q7W980
C	-24	SER	-	EXPRESSION TAG	UNP Q7W980
C	-23	ASN	-	EXPRESSION TAG	UNP Q7W980
C	-22	ALA	-	EXPRESSION TAG	UNP Q7W980
D	-24	SER	-	EXPRESSION TAG	UNP Q7W980
D	-23	ASN	-	EXPRESSION TAG	UNP Q7W980
D	-22	ALA	-	EXPRESSION TAG	UNP Q7W980

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



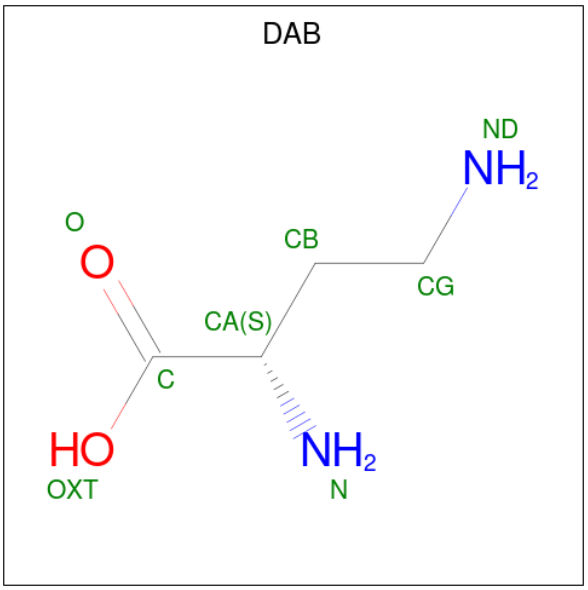
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	D	1	5	4	1	0	0
2	D	1	5	4	1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	B	1	6	3	3	0	0
			Total	C	O		
3	C	1	6	3	3	0	0

- Molecule 4 is 2,4-DIAMINOBUTYRIC ACID (three-letter code: DAB) (formula: $C_4H_{10}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	D	1	8	4	2	2	0	0
			Total	C	N	O		
4	D	1	8	4	2	2	0	0

- Molecule 5 is water.

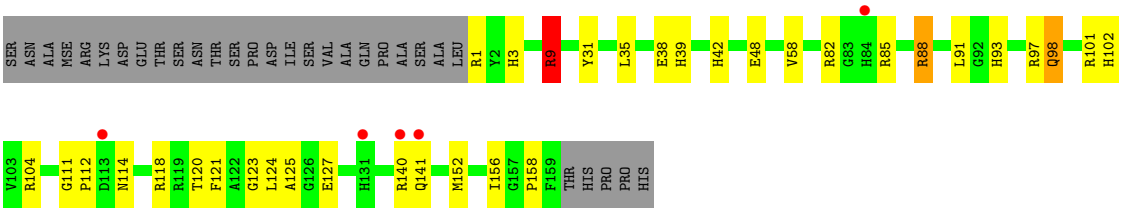
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	97	97	97	0	0
			Total	O		
5	B	91	91	91	0	0
			Total	O		
5	C	101	101	101	0	0
			Total	O		
5	D	88	88	88	0	0

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.

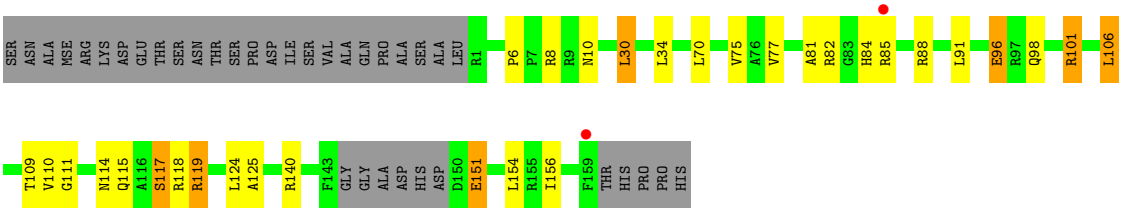
• Molecule 1: L-2,4-diaminobutyric acid acetyltransferase

Chain A:



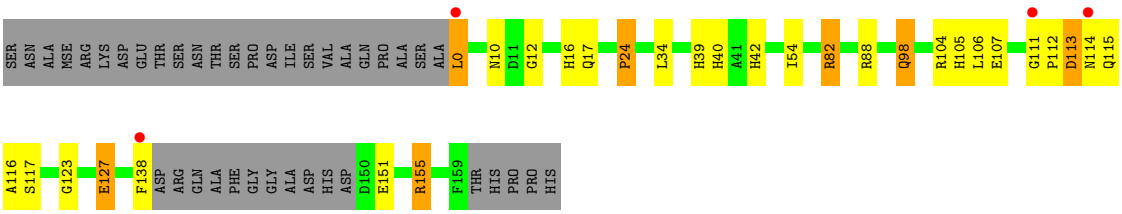
• Molecule 1: L-2,4-diaminobutyric acid acetyltransferase

Chain B:



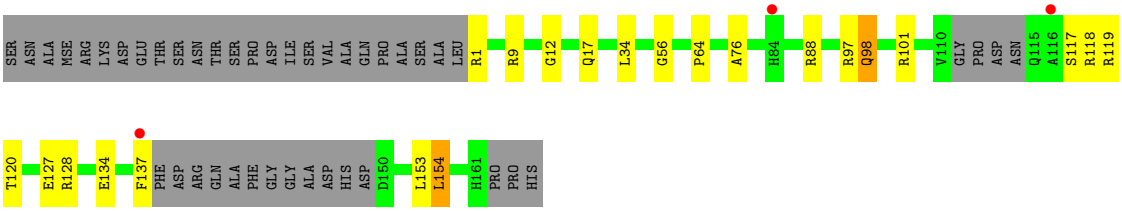
• Molecule 1: L-2,4-diaminobutyric acid acetyltransferase

Chain C:



• Molecule 1: L-2,4-diaminobutyric acid acetyltransferase

Chain D:


SER
ASN
ALA
MSE
ARG
LYS
ASP
GLU
THR
SER
ASN
THR
SER
PRO
ASP
ILE
SER
VAL
ALA
GLN
PRO
ALA
SER
ALA
LEU
R1
R9
G12
Q17
L34
G56
P64
A76
H84
R88
R97
Q98
R101
V110
GLY
PRO
ASP
ASN
Q115
A116
S117
R118
R119

T120
E127
R128
E134
F137
PHE
ASP
ARG
GLN
ALA
PHE
GLY
GLY
ALA
ASP
HIS
ASP
D150
L153
L154
H161
PRO
PRO
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	69.16Å 63.15Å 77.33Å 90.00° 96.06° 90.00°	Depositor
Resolution (Å)	48.80 – 1.87 48.80 – 1.86	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.80-1.87) 99.5 (48.80-1.86)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.61 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.165 , 0.210 0.165 , 0.211	Depositor DCC
R_{free} test set	2802 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	29.1	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 54.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	5650	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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: SO4, DAB, GOL

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	0.90	1/1383 (0.1%)	0.88	1/1869 (0.1%)
1	B	0.90	0/1365	0.87	1/1841 (0.1%)
1	C	0.95	0/1288	0.92	3/1737 (0.2%)
1	D	0.89	0/1261	0.86	1/1701 (0.1%)
All	All	0.91	1/5297 (0.0%)	0.88	6/7148 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	58	VAL	CB-CG1	5.95	1.65	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	9	ARG	NE-CZ-NH1	6.65	123.62	120.30
1	C	155[A]	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	C	155[B]	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	B	34	LEU	CB-CG-CD1	-5.71	101.29	111.00
1	D	97	ARG	NE-CZ-NH1	-5.59	117.51	120.30
1	C	82	ARG	NE-CZ-NH2	-5.39	117.60	120.30

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	1349	0	1290	36	0
1	B	1333	0	1293	26	0
1	C	1259	0	1226	37	0
1	D	1232	0	1204	26	0
2	A	15	0	0	0	0
2	B	20	0	0	2	0
2	C	15	0	0	1	0
2	D	10	0	0	3	0
3	A	12	0	16	0	0
3	B	6	0	8	2	0
3	C	6	0	8	2	0
4	D	16	0	18	1	0
5	A	97	0	0	8	0
5	B	91	0	0	3	0
5	C	101	0	0	8	0
5	D	88	0	0	8	0
All	All	5650	0	5063	116	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 (116) 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:D:9[B]:ARG:HD2	5:D:254:HOH:O	1.58	1.04
1:D:134:GLU:HG2	1:D:154:LEU:HD12	1.43	1.00
1:A:152:MSE:HG2	5:A:226:HOH:O	1.60	1.00
1:C:138[A]:PHE:CE2	1:D:9[A]:ARG:HG3	1.95	1.00
1:A:88:ARG:HD3	5:A:179:HOH:O	1.62	1.00
1:C:138[A]:PHE:CD2	1:D:9[A]:ARG:NH1	2.32	0.96
1:C:105:HIS:ND1	5:C:335:HOH:O	2.02	0.91
1:A:9:ARG:HG3	1:A:38:GLU:OE1	1.69	0.90
1:B:114:ASN:O	1:B:117:SER:HB2	1.76	0.84
1:A:9:ARG:CG	1:A:38:GLU:OE1	2.25	0.83
1:C:98[A]:GLN:H	1:C:98[A]:GLN:HE21	1.26	0.82
1:C:10:ASN:HB2	5:C:308:HOH:O	1.78	0.81
1:A:88:ARG:HD2	1:A:120:THR:HG23	1.61	0.80
1:C:138[A]:PHE:HE2	1:D:9[A]:ARG:HG3	1.42	0.80
1:C:138[A]:PHE:CE2	1:D:9[A]:ARG:NH1	2.49	0.79
1:C:114[A]:ASN:O	1:C:117:SER:HB2	1.82	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:ARG:HD3	2:C:166:SO4:O2	1.83	0.78
1:D:134:GLU:HG2	1:D:154:LEU:CD1	2.12	0.77
1:C:10:ASN:CB	5:C:308:HOH:O	2.34	0.76
1:B:98:GLN:O	1:B:101:ARG:HG2	1.88	0.74
1:B:111:GLY:HA3	1:B:151[A]:GLU:HG2	1.70	0.73
1:A:112:PRO:HA	1:A:152:MSE:HE1	1.73	0.71
1:B:119[B]:ARG:HG2	5:B:286:HOH:O	1.90	0.71
1:C:123:GLY:O	1:C:127:GLU:HG2	1.92	0.70
1:A:158:PRO:HD3	5:A:267:HOH:O	1.92	0.69
1:A:121:PHE:HB3	1:A:156:ILE:HD11	1.75	0.69
1:C:39:HIS:O	1:C:42[A]:HIS:HE1	1.76	0.69
1:A:118:ARG:HH22	1:A:152:MSE:CE	2.04	0.68
1:A:88:ARG:CD	5:A:179:HOH:O	2.32	0.68
1:D:9[A]:ARG:HB3	5:D:286:HOH:O	1.94	0.68
1:C:114[B]:ASN:O	1:C:117:SER:HB2	1.93	0.68
1:D:127[B]:GLU:OE2	5:D:298:HOH:O	2.11	0.67
1:A:3:HIS:HE1	1:A:48:GLU:OE1	1.78	0.67
4:D:168:DAB:HG3	5:D:245:HOH:O	1.95	0.66
1:C:138[A]:PHE:CZ	1:D:9[A]:ARG:HG3	2.32	0.63
1:D:120:THR:OG1	2:D:166:SO4:O1	2.04	0.63
1:C:138[A]:PHE:HE2	1:D:9[A]:ARG:CG	2.10	0.62
1:B:110:VAL:HG22	1:B:154:LEU:CD1	2.30	0.61
1:A:98[A]:GLN:O	1:A:101[A]:ARG:HD3	2.01	0.60
1:D:88:ARG:HD3	5:D:298:HOH:O	2.00	0.60
1:C:155[B]:ARG:HG2	5:C:335:HOH:O	2.01	0.59
1:B:110:VAL:HG22	1:B:154:LEU:HD13	1.84	0.58
1:B:8[B]:ARG:HD3	3:B:169:GOL:C2	2.33	0.58
1:C:138[A]:PHE:CE2	1:D:9[A]:ARG:CG	2.82	0.57
1:C:82:ARG:CG	1:C:82:ARG:HH11	2.17	0.56
1:A:118:ARG:NH2	1:A:152:MSE:CE	2.69	0.56
1:B:81:ALA:O	1:B:84:HIS:HD2	1.89	0.55
1:C:104:ARG:HG2	1:C:105:HIS:CD2	2.42	0.54
1:D:88:ARG:CD	5:D:298:HOH:O	2.55	0.54
1:D:88:ARG:NH2	2:D:166:SO4:O2	2.39	0.54
1:A:118:ARG:HH22	1:A:152:MSE:HE2	1.71	0.54
1:C:24:PRO:HG2	1:C:113:ASP:O	2.08	0.54
1:B:85:ARG:HA	2:B:165:SO4:O1	2.09	0.53
1:D:17[B]:GLN:OE1	5:D:237:HOH:O	2.18	0.53
1:A:118:ARG:NH2	1:A:152:MSE:HE2	2.24	0.52
1:A:125:ALA:HB2	1:A:156:ILE:HG23	1.92	0.52
1:D:127[B]:GLU:OE1	1:D:128[B]:ARG:NH1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98[A]:GLN:CD	1:A:98[A]:GLN:H	2.13	0.52
1:B:75[B]:VAL:HG23	1:B:75[B]:VAL:O	2.10	0.51
1:A:101[B]:ARG:HE	1:A:102:HIS:CE1	2.27	0.51
1:C:82:ARG:HH11	1:C:82:ARG:HG2	1.75	0.51
1:C:54:ILE:O	3:C:168:GOL:H11	2.11	0.51
1:C:114[B]:ASN:ND2	1:C:117:SER:OG	2.44	0.50
1:B:8[B]:ARG:HD3	3:B:169:GOL:H2	1.92	0.50
1:C:82:ARG:NH2	5:C:289:HOH:O	2.42	0.50
1:B:6:PRO:HD3	2:B:168:SO4:O2	2.11	0.50
1:D:12:GLY:HA3	1:D:34:LEU:HD23	1.94	0.50
1:C:12:GLY:HA3	1:C:34:LEU:HD23	1.93	0.49
1:A:112:PRO:CA	1:A:152:MSE:HE1	2.41	0.49
1:A:9:ARG:CD	1:A:38:GLU:OE1	2.59	0.49
1:A:111:GLY:CA	1:A:114:ASN:HD22	2.24	0.49
1:B:98:GLN:O	1:B:101:ARG:CG	2.60	0.49
1:A:31:TYR:OH	1:B:109:THR:OG1	2.25	0.48
1:B:70:LEU:HB3	1:B:106:LEU:CD2	2.44	0.48
1:B:96[A]:GLU:HG3	5:B:278:HOH:O	2.13	0.48
1:A:82:ARG:HA	1:A:82:ARG:HD3	1.60	0.48
1:A:9:ARG:HD3	1:A:38:GLU:OE1	2.14	0.47
1:B:85:ARG:HD2	1:B:88:ARG:CZ	2.44	0.47
1:B:115:GLN:HE22	1:B:118:ARG:HE	1.63	0.46
1:D:88:ARG:HB2	2:D:166:SO4:O4	2.15	0.46
1:A:31:TYR:CE1	1:A:35:LEU:HD11	2.51	0.46
1:A:141[A]:GLN:NE2	1:B:10:ASN:HD21	2.14	0.45
1:A:88:ARG:HG2	5:A:179:HOH:O	2.15	0.45
1:B:82:ARG:NH2	5:B:316:HOH:O	2.48	0.45
1:B:110:VAL:CG2	1:B:154:LEU:HD11	2.47	0.45
1:A:140[A]:ARG:HG3	5:A:209:HOH:O	2.17	0.44
1:A:123:GLY:O	1:A:127[B]:GLU:HG3	2.18	0.44
1:D:137:PHE:HB2	1:D:153:LEU:HB2	1.99	0.44
1:B:125:ALA:HB2	1:B:156:ILE:HG23	2.00	0.43
1:B:91:LEU:HB3	1:B:124:LEU:HD22	2.00	0.43
1:A:104:ARG:O	1:A:158:PRO:HA	2.19	0.43
1:B:110:VAL:HG22	1:B:154:LEU:HD11	2.01	0.43
1:A:91:LEU:HB3	1:A:124:LEU:HD21	2.00	0.43
1:C:111:GLY:HA3	1:C:151[B]:GLU:OE2	2.19	0.43
1:C:116:ALA:HA	5:C:250:HOH:O	2.19	0.43
1:C:0:LEU:HD23	1:C:0:LEU:HA	1.84	0.43
1:B:30:LEU:HD12	1:B:30:LEU:O	2.19	0.42
1:C:111:GLY:O	1:C:112:PRO:C	2.56	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ASN:ND2	5:A:261:HOH:O	2.51	0.42
1:A:3:HIS:CE1	1:A:48:GLU:HB3	2.55	0.42
1:D:117:SER:C	1:D:119:ARG:N	2.72	0.42
1:D:9[B]:ARG:CD	5:D:254:HOH:O	2.37	0.42
1:C:98[A]:GLN:H	1:C:98[A]:GLN:NE2	2.05	0.41
1:D:56:GLY:HA2	1:D:76:ALA:O	2.19	0.41
1:A:39:HIS:O	1:A:42[B]:HIS:HE1	2.02	0.41
1:C:16:HIS:HE1	5:C:280:HOH:O	2.03	0.41
1:C:40:HIS:HA	1:C:42[A]:HIS:CE1	2.55	0.41
1:C:17:GLN:HG2	3:C:168:GOL:H32	2.02	0.41
1:D:98:GLN:HE21	1:D:98:GLN:HB2	1.68	0.41
1:A:93:HIS:CE1	1:A:97:ARG:HH21	2.38	0.41
1:B:106:LEU:HB2	1:B:156:ILE:HB	2.03	0.41
1:C:82:ARG:NH1	5:C:283:HOH:O	2.52	0.41
1:C:106:LEU:HD23	1:C:107:GLU:N	2.35	0.41
1:C:42[A]:HIS:HE2	1:D:64:PRO:HB2	1.86	0.40
1:A:88:ARG:CG	5:A:179:HOH:O	2.66	0.40
1:C:106:LEU:HD23	1:C:106:LEU:C	2.41	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	167/189 (88%)	164 (98%)	3 (2%)	0	100	100
1	B	161/189 (85%)	159 (99%)	2 (1%)	0	100	100
1	C	152/189 (80%)	149 (98%)	3 (2%)	0	100	100
1	D	147/189 (78%)	146 (99%)	1 (1%)	0	100	100
All	All	627/756 (83%)	618 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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	140/153 (92%)	134 (96%)	6 (4%)	29	17
1	B	139/153 (91%)	127 (91%)	12 (9%)	10	3
1	C	132/153 (86%)	125 (95%)	7 (5%)	22	11
1	D	129/153 (84%)	124 (96%)	5 (4%)	32	20
All	All	540/612 (88%)	510 (94%)	30 (6%)	24	10

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ARG
1	A	9	ARG
1	A	85	ARG
1	A	88	ARG
1	A	98[A]	GLN
1	A	98[B]	GLN
1	B	30	LEU
1	B	77	VAL
1	B	96[A]	GLU
1	B	96[B]	GLU
1	B	101	ARG
1	B	106	LEU
1	B	117	SER
1	B	119[A]	ARG
1	B	119[B]	ARG
1	B	140	ARG
1	B	151[A]	GLU
1	B	151[B]	GLU
1	C	0	LEU
1	C	24	PRO
1	C	98[A]	GLN
1	C	98[B]	GLN
1	C	113	ASP
1	C	115	GLN

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Mol	Chain	Res	Type
1	C	127	GLU
1	D	1	ARG
1	D	98	GLN
1	D	101	ARG
1	D	118	ARG
1	D	154	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	HIS
1	A	114	ASN
1	B	84	HIS
1	B	98	GLN
1	B	115	GLN
1	C	16	HIS
1	D	98	GLN
1	D	102	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

18 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	SO4	C	167	-	4,4,4	0.26	0	6,6,6	0.21	0
4	DAB	D	168	-	6,7,7	1.16	0	4,8,8	2.73	1 (25%)
2	SO4	B	166	-	4,4,4	0.29	0	6,6,6	0.32	0
2	SO4	D	166	-	4,4,4	0.21	0	6,6,6	0.39	0
3	GOL	C	168	-	5,5,5	0.66	0	5,5,5	0.79	0
4	DAB	D	167	-	6,7,7	1.23	1 (16%)	4,8,8	3.65	3 (75%)
2	SO4	A	167	-	4,4,4	0.35	0	6,6,6	0.59	0
2	SO4	B	165	-	4,4,4	0.26	0	6,6,6	0.45	0
2	SO4	A	165	-	4,4,4	0.35	0	6,6,6	0.31	0
2	SO4	C	166	-	4,4,4	0.34	0	6,6,6	0.76	0
2	SO4	B	168	-	4,4,4	0.32	0	6,6,6	0.47	0
3	GOL	A	169	-	5,5,5	0.46	0	5,5,5	1.52	0
2	SO4	D	165	-	4,4,4	0.25	0	6,6,6	0.56	0
3	GOL	B	169	-	5,5,5	0.48	0	5,5,5	1.82	1 (20%)
2	SO4	C	165	-	4,4,4	0.57	0	6,6,6	0.72	0
2	SO4	A	166	-	4,4,4	0.38	0	6,6,6	0.64	0
2	SO4	B	167	-	4,4,4	0.27	0	6,6,6	0.45	0
3	GOL	A	168	-	5,5,5	0.52	0	5,5,5	0.99	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
4	DAB	D	168	-	-	4/7/7/7	-
3	GOL	C	168	-	-	2/4/4/4	-
4	DAB	D	167	-	-	3/7/7/7	-
3	GOL	B	169	-	-	3/4/4/4	-
3	GOL	A	169	-	-	4/4/4/4	-
3	GOL	A	168	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	167	DAB	O-C	2.51	1.29	1.22

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	D	168	DAB	OXT-C-O	-5.41	111.80	124.08
4	D	167	DAB	OXT-C-O	-5.36	111.92	124.08
4	D	167	DAB	CB-CA-N	-3.65	100.59	110.12
4	D	167	DAB	CB-CA-C	3.30	119.19	110.45
3	B	169	GOL	O2-C2-C1	-3.05	96.57	109.18

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	168	GOL	O1-C1-C2-C3
3	A	169	GOL	O1-C1-C2-C3
3	A	169	GOL	C1-C2-C3-O3
3	B	169	GOL	C1-C2-C3-O3
3	C	168	GOL	C1-C2-C3-O3
4	D	168	DAB	N-CA-CB-CG
4	D	168	DAB	C-CA-CB-CG
3	B	169	GOL	O1-C1-C2-C3
3	A	168	GOL	O1-C1-C2-O2
3	A	169	GOL	O1-C1-C2-O2
3	B	169	GOL	O2-C2-C3-O3
3	A	169	GOL	O2-C2-C3-O3
3	C	168	GOL	O2-C2-C3-O3
4	D	168	DAB	O-C-CA-CB
4	D	168	DAB	OXT-C-CA-CB
4	D	167	DAB	O-C-CA-CB
4	D	167	DAB	OXT-C-CA-CB
4	D	167	DAB	CA-CB-CG-ND

There are no ring outliers.

7 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	168	DAB	1	0
2	D	166	SO4	3	0
3	C	168	GOL	2	0
2	B	165	SO4	1	0
2	C	166	SO4	1	0
2	B	168	SO4	1	0
3	B	169	GOL	2	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 (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	157/189 (83%)	0.14	5 (3%)	47	49	25, 33, 46, 54	0
1	B	151/189 (79%)	0.19	2 (1%)	77	79	25, 34, 50, 64	1 (0%)
1	C	147/189 (77%)	0.15	4 (2%)	54	56	23, 30, 49, 59	0
1	D	143/189 (75%)	0.04	3 (2%)	63	65	25, 33, 55, 73	0
All	All	598/756 (79%)	0.13	14 (2%)	60	62	23, 32, 50, 73	1 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	116	ALA	4.6
1	C	111	GLY	4.5
1	C	138[A]	PHE	4.2
1	D	84	HIS	3.3
1	A	113	ASP	3.1
1	A	140[A]	ARG	2.9
1	A	131[A]	HIS	2.7
1	C	0	LEU	2.7
1	B	159	PHE	2.5
1	B	85	ARG	2.4
1	A	141[A]	GLN	2.2
1	D	137	PHE	2.1
1	C	114[A]	ASN	2.1
1	A	84	HIS	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 monosaccharides 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(Å^2)	Q<0.9
2	SO4	C	167	5/5	0.77	0.23	111,112,112,113	0
4	DAB	D	167	8/8	0.81	0.26	27,36,39,40	0
3	GOL	B	169	6/6	0.86	0.31	29,38,41,42	0
2	SO4	B	168	5/5	0.86	0.34	82,82,82,84	0
4	DAB	D	168	8/8	0.86	0.29	29,36,37,39	0
3	GOL	A	168	6/6	0.90	0.20	53,57,58,58	0
2	SO4	B	166	5/5	0.90	0.11	83,84,84,85	0
3	GOL	C	168	6/6	0.91	0.23	46,50,52,52	0
2	SO4	B	167	5/5	0.93	0.15	79,80,81,81	0
2	SO4	D	165	5/5	0.93	0.10	62,63,64,67	0
2	SO4	A	165	5/5	0.93	0.15	64,64,68,68	0
3	GOL	A	169	6/6	0.93	0.15	35,35,36,38	0
2	SO4	D	166	5/5	0.95	0.24	63,67,67,70	0
2	SO4	B	165	5/5	0.96	0.17	62,63,64,67	0
2	SO4	A	167	5/5	0.97	0.10	58,59,60,60	0
2	SO4	C	166	5/5	0.98	0.07	40,41,47,48	0
2	SO4	C	165	5/5	0.98	0.08	40,40,43,43	0
2	SO4	A	166	5/5	0.99	0.07	48,50,51,51	0

6.5 Other polymers (i)

There are no such residues in this entry.