



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 12:31 AM GMT

PDB ID : 2R5S
Title : The crystal structure of a domain of protein VP0806 (unknown function) from *Vibrio parahaemolyticus* RIMD 2210633
Authors : Tan, K.; Wu, R.; Abdullah, J.; Freeman, L.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2007-09-04
Resolution : 2.14 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

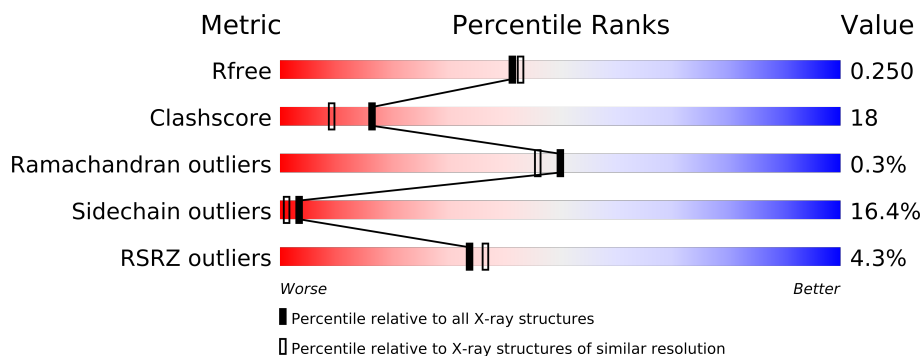
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.14 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}	66092	1116 (2.16-2.12)
Clashscore	79885	1302 (2.16-2.12)
Ramachandran outliers	78287	1281 (2.16-2.12)
Sidechain outliers	78261	1281 (2.16-2.12)
RSRZ outliers	66119	1116 (2.16-2.12)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	176	
1	B	176	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	MES	B	2	-	X
3	MES	B	285	-	X
4	EDO	A	2	-	X
4	EDO	A	3	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 2850 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

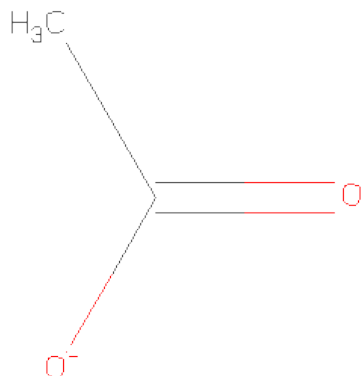
- Molecule 1 is a protein called Uncharacterized protein VP0806.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	Se	0	2	0
			1370	864	229	274	2	1			
1	B	171	Total	C	N	O	S	Se	0	0	0
			1356	853	228	272	2	1			

There are 6 discrepancies between the modelled and reference sequences:

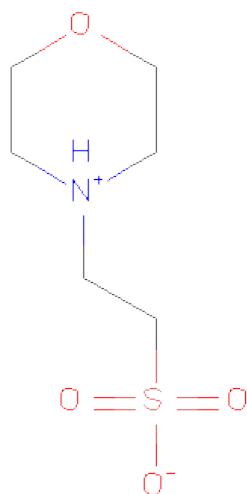
Chain	Residue	Modelled	Actual	Comment	Reference
A	109	SER	-	EXPRESSION TAG	UNP Q87RI8
A	110	ASN	-	EXPRESSION TAG	UNP Q87RI8
A	111	ALA	-	EXPRESSION TAG	UNP Q87RI8
B	109	SER	-	EXPRESSION TAG	UNP Q87RI8
B	110	ASN	-	EXPRESSION TAG	UNP Q87RI8
B	111	ALA	-	EXPRESSION TAG	UNP Q87RI8

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONICACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 5 is water.

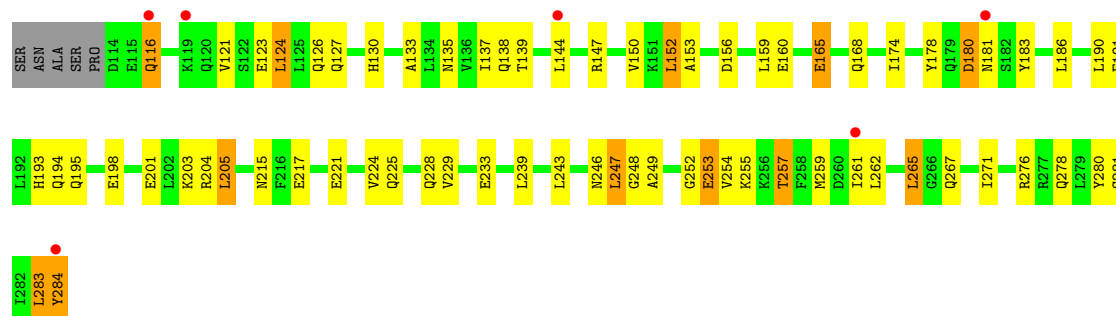
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	31	Total	O	0	0
			31	31		
5	B	29	Total	O	0	0
			29	29		

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.

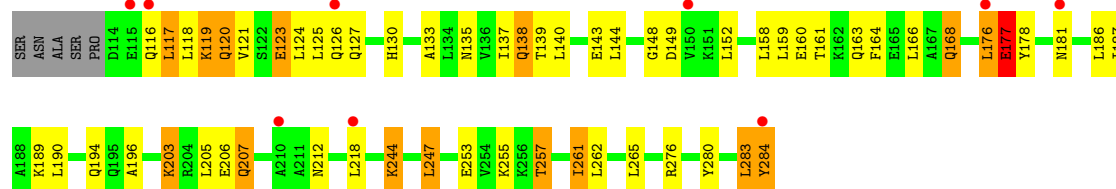
- Molecule 1: Uncharacterized protein VP0806

Chain A: 



- Molecule 1: Uncharacterized protein VP0806

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	53.00Å 116.48Å 62.22Å 90.00° 88.75° 90.00°	Depositor
Resolution (Å)	29.12 – 2.14 29.12 – 2.14	Depositor EDS
% Data completeness (in resolution range)	95.0 (29.12-2.14) 90.0 (29.12-2.14)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.223 , 0.279 0.215 , 0.250	Depositor DCC
R_{free} test set	1011 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.574	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 85.5	EDS
Estimated twinning fraction	0.329 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 19739 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2850	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, MES, ACT

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	0.97	2/1390 (0.1%)	0.99	1/1874 (0.1%)
1	B	1.05	1/1370 (0.1%)	1.06	2/1848 (0.1%)
All	All	1.01	3/2760 (0.1%)	1.02	3/3722 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	165	GLU	CB-CG	-5.53	1.41	1.52
1	A	253	GLU	CG-CD	5.24	1.59	1.51
1	B	177	GLU	CG-CD	-5.21	1.44	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	244	LYS	CD-CE-NZ	-5.47	99.11	111.70
1	A	124	LEU	CA-CB-CG	5.47	127.89	115.30
1	B	176	LEU	CA-CB-CG	5.32	127.55	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1370	0	1374	52	1
1	B	1356	0	1359	52	1
2	B	4	0	3	1	0
3	A	12	0	12	1	0
3	B	36	0	36	0	0
4	A	8	0	12	7	0
4	B	4	0	6	0	0
5	A	31	0	0	4	0
5	B	29	0	0	1	0
All	All	2850	0	2802	101	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 18.

All (101) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:161:THR:HG23	1:B:163:GLN:HG3	1.39	1.00
1:A:281:SER:HB3	4:A:2:EDO:H21	1.46	0.98
1:A:224:VAL:O	1:A:228:GLN:HG3	1.67	0.94
1:A:168:GLN:HB2	1:A:190:LEU:HD21	1.46	0.93
1:A:123:GLU:O	1:A:127:GLN:HG3	1.74	0.87
1:A:252:GLY:HA2	5:A:288:HOH:O	1.78	0.83
1:A:281:SER:CB	4:A:2:EDO:H21	2.08	0.83
1:B:158:LEU:O	1:B:161:THR:HG22	1.79	0.81
1:A:281:SER:HB3	4:A:2:EDO:C2	2.15	0.77
1:B:168:GLN:HG2	1:B:190:LEU:HD21	1.64	0.77
1:B:117:LEU:HD23	1:B:117:LEU:H	1.47	0.77
1:B:133:ALA:O	1:B:137:ILE:HG12	1.87	0.74
1:A:243:LEU:HD21	1:A:254:VAL:HG12	1.71	0.72
1:A:204:ARG:NH2	1:A:205:LEU:HD13	2.04	0.72
1:B:161:THR:CG2	1:B:163:GLN:HE21	2.04	0.70
1:B:207:GLN:HA	1:B:207:GLN:HE21	1.56	0.70
1:A:168:GLN:CB	1:A:190:LEU:HD21	2.23	0.68
1:B:164:PHE:O	1:B:168:GLN:HB2	1.94	0.68
1:A:278:GLN:HA	5:A:315:HOH:O	1.94	0.67
1:B:190:LEU:O	1:B:194:GLN:HG3	1.95	0.67
1:A:217:GLU:O	1:A:221:GLU:HG3	1.94	0.66
1:A:283:LEU:HD23	1:B:189:LYS:HG3	1.78	0.66
1:A:133:ALA:O	1:A:137:ILE:HG12	1.97	0.65
1:A:152:LEU:HD13	1:A:174:ILE:HD12	1.79	0.64
1:A:281:SER:CA	4:A:2:EDO:H21	2.28	0.64
1:B:121:VAL:HG11	1:B:137:ILE:HD11	1.78	0.64
1:A:247:LEU:HD23	1:A:284[B]:TYR:CE2	2.32	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:130:HIS:CE1	1:A:160:GLU:HG3	2.34	0.62
1:A:257:THR:O	1:A:261:ILE:HD13	2.00	0.61
1:B:280:TYR:HA	1:B:283:LEU:HD22	1.82	0.61
1:A:243:LEU:CD2	1:A:254:VAL:HG12	2.31	0.60
1:B:203:LYS:NZ	1:B:206:GLU:OE2	2.36	0.58
1:B:161:THR:CG2	1:B:163:GLN:HG3	2.25	0.57
1:A:280:TYR:HA	1:A:283:LEU:HD22	1.86	0.57
1:A:229:VAL:HG12	4:A:3:EDO:H22	1.87	0.57
1:B:203:LYS:O	1:B:207:GLN:HG2	2.03	0.56
1:A:135:ASN:O	1:A:139:THR:HG23	2.05	0.56
1:A:262:LEU:HD13	1:A:276:ARG:HG2	1.87	0.56
1:A:204:ARG:NH2	1:A:205:LEU:CD1	2.68	0.56
1:B:161:THR:HG21	1:B:163:GLN:HE21	1.71	0.56
1:A:284[A]:TYR:HD2	1:B:181:ASN:ND2	2.04	0.55
1:A:253:GLU:O	1:A:257:THR:HG22	2.07	0.55
1:B:161:THR:HG23	1:B:163:GLN:HE21	1.70	0.54
1:A:191:GLU:O	1:A:195:GLN:HG3	2.07	0.54
1:B:161:THR:HG23	1:B:163:GLN:CG	2.25	0.54
1:B:253:GLU:O	1:B:257:THR:CG2	2.57	0.53
1:B:161:THR:CG2	1:B:163:GLN:NE2	2.73	0.51
1:B:207:GLN:CA	1:B:207:GLN:HE21	2.23	0.51
1:B:135:ASN:O	1:B:139:THR:HG23	2.12	0.49
1:B:244:LYS:CE	2:B:1:ACT:H2	2.43	0.49
1:A:281:SER:HA	4:A:2:EDO:H21	1.94	0.49
3:A:4:MES:H71	5:A:287:HOH:O	2.12	0.49
1:B:161:THR:HG21	1:B:163:GLN:NE2	2.26	0.49
1:A:284[A]:TYR:CD2	1:B:181:ASN:ND2	2.80	0.49
1:A:253:GLU:O	1:A:257:THR:CG2	2.60	0.49
1:A:247:LEU:CD2	1:A:284[B]:TYR:CE2	2.97	0.47
1:A:267:GLN:NE2	1:B:196:ALA:HA	2.29	0.47
1:A:248:GLY:O	1:A:249:ALA:C	2.51	0.47
1:B:117:LEU:CD2	1:B:117:LEU:H	2.23	0.47
1:B:116:GLN:NE2	1:B:120:GLN:HE22	2.12	0.47
1:B:125:LEU:O	1:B:126:GLN:C	2.51	0.47
1:A:121:VAL:HG21	1:A:137:ILE:CD1	2.45	0.46
1:B:253:GLU:O	1:B:257:THR:HG23	2.15	0.46
1:A:180:ASP:HB3	1:A:183:TYR:H	1.78	0.46
1:A:243:LEU:HD21	1:A:255:LYS:N	2.31	0.46
1:B:177:GLU:HG3	1:B:177:GLU:H	1.50	0.46
1:B:116:GLN:HG2	1:B:116:GLN:O	2.14	0.46
1:A:193:HIS:ND1	4:A:3:EDO:H12	2.31	0.46
1:A:233:GLU:HG3	1:A:271:ILE:HD12	1.98	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:116:GLN:HG2	1:B:119:LYS:HB3	1.99	0.45
1:B:117:LEU:HA	1:B:120:GLN:HB2	1.98	0.45
1:B:148:GLY:HA3	1:B:178:TYR:CD2	2.52	0.45
1:A:121:VAL:HG21	1:A:137:ILE:HD13	1.98	0.44
1:B:121:VAL:HG21	1:B:137:ILE:HD13	1.99	0.44
1:B:121:VAL:CG1	1:B:137:ILE:HD11	2.47	0.44
1:B:253:GLU:O	1:B:257:THR:HG22	2.18	0.43
1:A:201:GLU:OE2	1:A:204:ARG:NH2	2.42	0.43
1:B:123:GLU:O	1:B:127:GLN:HG3	2.17	0.43
1:B:247:LEU:HD22	5:B:294:HOH:O	2.17	0.43
1:B:261:ILE:CD1	1:B:261:ILE:N	2.81	0.43
1:A:225:GLN:OE1	1:A:225:GLN:HA	2.19	0.43
1:B:138:GLN:HB2	1:B:138:GLN:HE21	1.61	0.43
1:B:284:TYR:C	1:B:284:TYR:HD2	2.21	0.43
1:A:255:LYS:O	1:A:259:MSE:HG2	2.18	0.43
1:B:149:ASP:OD1	1:B:149:ASP:N	2.52	0.43
1:B:118:LEU:HG	1:B:144:LEU:HD21	2.01	0.43
1:B:130:HIS:CE1	1:B:160:GLU:HG3	2.55	0.42
1:B:284:TYR:C	1:B:284:TYR:CD2	2.93	0.42
1:B:262:LEU:HD13	1:B:276:ARG:HG2	2.02	0.42
1:A:156:ASP:O	1:A:160:GLU:HG2	2.20	0.42
1:B:118:LEU:HD11	1:B:144:LEU:HD11	2.00	0.42
1:A:265:LEU:HA	1:A:265:LEU:HD12	1.78	0.42
1:A:150:VAL:O	1:A:153:ALA:HB3	2.20	0.41
1:B:262:LEU:HD13	1:B:276:ARG:CG	2.50	0.41
1:B:257:THR:O	1:B:261:ILE:HD13	2.21	0.41
1:A:144:LEU:HD12	1:A:147:ARG:HD3	2.03	0.40
1:A:190:LEU:O	1:A:194:GLN:HG2	2.22	0.40
1:A:116:GLN:N	5:A:301:HOH:O	2.45	0.40
1:A:123:GLU:O	1:A:127:GLN:CG	2.58	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:178:TYR:OH	1:B:178:TYR:OH[4.556]	1.87	0.33

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	170/176 (97%)	160 (94%)	9 (5%)	1 (1%)	33	25
1	B	169/176 (96%)	164 (97%)	5 (3%)	0	100	100
All	All	339/352 (96%)	324 (96%)	14 (4%)	1 (0%)	50	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	215	ASN

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	148/149 (99%)	126 (85%)	22 (15%)	4	1
1	B	146/149 (98%)	118 (81%)	28 (19%)	2	1
All	All	294/298 (99%)	244 (83%)	50 (17%)	3	1

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	124	LEU
1	A	126	GLN
1	A	138	GLN
1	A	152	LEU
1	A	159	LEU
1	A	165	GLU
1	A	180	ASP

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Mol	Chain	Res	Type
1	A	181[A]	ASN
1	A	181[B]	ASN
1	A	186	LEU
1	A	198	GLU
1	A	203	LYS
1	A	205	LEU
1	A	239	LEU
1	A	246	ASN
1	A	247	LEU
1	A	257	THR
1	A	265	LEU
1	A	283	LEU
1	A	284[A]	TYR
1	A	284[B]	TYR
1	B	117	LEU
1	B	119	LYS
1	B	120	GLN
1	B	123	GLU
1	B	124	LEU
1	B	138	GLN
1	B	140	LEU
1	B	143	GLU
1	B	152	LEU
1	B	159	LEU
1	B	166	LEU
1	B	168	GLN
1	B	176	LEU
1	B	177	GLU
1	B	186	LEU
1	B	187	ILE
1	B	203	LYS
1	B	205	LEU
1	B	207	GLN
1	B	212	ASN
1	B	218	LEU
1	B	247	LEU
1	B	255	LYS
1	B	257	THR
1	B	261	ILE
1	B	265	LEU
1	B	283	LEU
1	B	284	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	130	HIS
1	A	246	ASN
1	A	267	GLN
1	B	116	GLN
1	B	130	HIS
1	B	138	GLN
1	B	163	GLN
1	B	195	GLN
1	B	207	GLN
1	B	241	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

8 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	$\# Z > 2$
4	EDO	A	2	-	3,3,3	0.59	0	2,2,2	0.10	0
4	EDO	A	3	-	3,3,3	0.66	0	2,2,2	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MES	A	4	-	12,12,12	2.06	1 (8%)	16,16,16	2.36	5 (31%)
2	ACT	B	1	-	1,3,3	2.47	1 (100%)	0,3,3	0.00	-
3	MES	B	2	-	12,12,12	1.78	2 (16%)	16,16,16	2.35	6 (37%)
3	MES	B	285	-	12,12,12	2.19	4 (33%)	16,16,16	2.98	8 (50%)
4	EDO	B	286	-	3,3,3	0.82	0	2,2,2	0.22	0
3	MES	B	3	-	12,12,12	1.85	1 (8%)	16,16,16	2.37	6 (37%)

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	EDO	A	2	-	-	0/1/1/1	0/0/0/0
4	EDO	A	3	-	-	0/1/1/1	0/0/0/0
3	MES	A	4	-	-	0/6/14/14	1/1/1/1
2	ACT	B	1	-	-	0/0/0/0	0/0/0/0
3	MES	B	2	-	-	0/6/14/14	0/1/1/1
3	MES	B	285	-	-	0/6/14/14	0/1/1/1
4	EDO	B	286	-	-	0/1/1/1	0/0/0/0
3	MES	B	3	-	-	0/6/14/14	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	4	MES	C8-S	-6.06	1.67	1.78
3	B	285	MES	C8-S	-5.77	1.68	1.78
3	B	3	MES	C8-S	-5.21	1.69	1.78
3	B	2	MES	C8-S	-4.60	1.70	1.78
2	B	1	ACT	CH3-C	2.47	1.52	1.48
3	B	285	MES	C7-N4	-2.24	1.44	1.50
3	B	2	MES	O3S-S	2.16	1.50	1.45
3	B	285	MES	O3S-S	2.13	1.50	1.45
3	B	285	MES	O2S-S	2.01	1.50	1.45

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	3	MES	C5-N4-C3	6.75	123.13	109.75
3	B	285	MES	O1-C6-C5	-5.40	104.91	111.34
3	B	285	MES	C5-N4-C3	5.39	120.44	109.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	4	MES	C5-N4-C3	5.31	120.27	109.75
3	B	2	MES	C5-N4-C3	5.25	120.16	109.75
3	B	285	MES	C6-C5-N4	-4.42	104.29	109.96
3	B	285	MES	O1-C2-C3	4.27	116.43	111.34
3	B	285	MES	C7-N4-C3	4.01	122.15	111.66
3	B	2	MES	O1-C6-C5	-3.95	106.64	111.34
3	A	4	MES	C7-N4-C3	3.56	120.97	111.66
3	A	4	MES	C6-C5-N4	-3.41	105.58	109.96
3	B	2	MES	C7-N4-C3	3.36	120.44	111.66
3	A	4	MES	C2-C3-N4	-3.27	105.76	109.96
3	B	2	MES	C2-C3-N4	-3.27	105.77	109.96
3	B	2	MES	C7-N4-C5	2.96	119.41	111.66
3	B	3	MES	C7-N4-C3	2.85	119.11	111.66
3	B	3	MES	C2-C3-N4	2.83	113.59	109.96
3	A	4	MES	C7-N4-C5	2.71	118.75	111.66
3	B	2	MES	O3S-S-C8	2.48	112.40	105.99
3	B	285	MES	O3S-S-C8	2.46	112.33	105.99
3	B	3	MES	C7-N4-C5	2.33	117.74	111.66
3	B	285	MES	O2S-S-O1S	-2.27	108.05	112.44
3	B	285	MES	C7-N4-C5	2.20	117.41	111.66
3	B	3	MES	C8-C7-N4	-2.19	108.69	112.44
3	B	3	MES	O3S-S-O1S	-2.15	106.94	112.48

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	4	MES	C2-C3-C5-C6-N4-O1

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 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	171/176 (97%)	0.63	6 (3%)	42	44	44, 56, 65, 74	1 (0%)
1	B	171/176 (97%)	0.66	9 (5%)	25	28	46, 56, 65, 73	1 (0%)
All	All	342/352 (97%)	0.65	15 (4%)	34	36	44, 56, 65, 74	2 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	284[A]	TYR	5.1
1	B	218	LEU	3.6
1	B	116	GLN	3.2
1	B	284	TYR	3.1
1	B	150	VAL	2.5
1	A	119	LYS	2.4
1	B	210	ALA	2.4
1	B	115	GLU	2.3
1	A	116	GLN	2.2
1	A	181[A]	ASN	2.1
1	A	261	ILE	2.1
1	B	126	GLN	2.1
1	B	176	LEU	2.1
1	A	144	LEU	2.0
1	B	181	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	EDO	A	3	4/4	0.27	6.75	66,68,68,69	0
4	EDO	A	2	4/4	0.28	6.30	59,59,62,63	0
3	MES	B	285	12/12	0.24	3.93	65,69,88,88	0
3	MES	B	2	12/12	0.23	2.90	64,67,85,86	12
3	MES	B	3	12/12	0.31	1.94	83,85,94,94	12
2	ACT	B	1	4/4	0.23	1.38	64,64,64,66	0
3	MES	A	4	12/12	0.25	1.08	75,79,93,93	12
4	EDO	B	286	4/4	0.22	0.65	72,73,74,75	0

6.5 Other polymers ⓘ

There are no such residues in this entry.