



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 04:30 pm BST

PDB ID : 3GFU
Title : FaeE-FaeG chaperone-major pilin complex of F4 ac 5/95 fimbriae
Authors : Van Molle, I.; Moonens, K.; Garcia-Pino, A.; Buts, L.; Bouckaert, J.; De Greve, H.
Deposited on : 2009-02-27
Resolution : 1.99 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

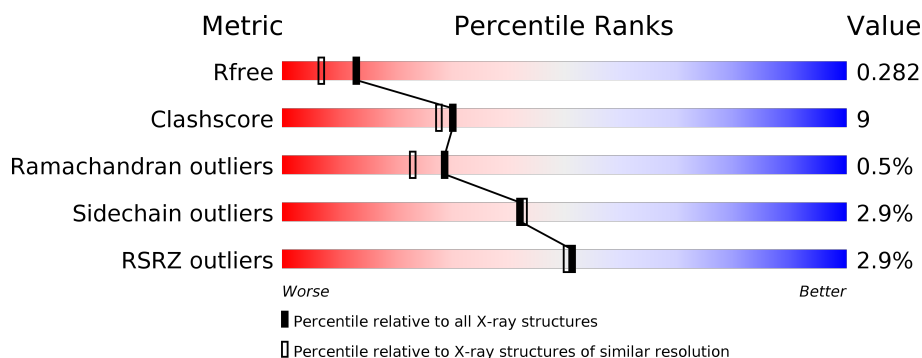
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	224	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>15%</div> <div>• 6%</div> </div> </div>
1	C	224	<div> <div>4%</div> <div> <div></div> <div>75%</div> <div>18%</div> <div>• 6%</div> </div> </div>
2	B	251	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>14%</div> <div>• 12%</div> </div> </div>
2	D	251	<div> <div>3%</div> <div> <div></div> <div>74%</div> <div>12%</div> <div>• 13%</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6775 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 Chaperone protein faeE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	210	Total	C	N	O	S	0	0	0
			1559	998	259	298	4			
1	A	211	Total	C	N	O	S	0	1	0
			1593	1016	272	301	4			

- Molecule 2 is a protein called FaeG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	219	Total	C	N	O	S	0	1	0
			1551	984	263	303	1			
2	B	220	Total	C	N	O	S	0	0	0
			1559	993	260	305	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	5	HIS	-	LINKER	UNP Q6T3W5
D	6	HIS	-	LINKER	UNP Q6T3W5
D	7	HIS	-	LINKER	UNP Q6T3W5
D	8	HIS	-	LINKER	UNP Q6T3W5
D	9	HIS	-	LINKER	UNP Q6T3W5
D	10	HIS	-	LINKER	UNP Q6T3W5
B	5	HIS	-	LINKER	UNP Q6T3W5
B	6	HIS	-	LINKER	UNP Q6T3W5
B	7	HIS	-	LINKER	UNP Q6T3W5
B	8	HIS	-	LINKER	UNP Q6T3W5
B	9	HIS	-	LINKER	UNP Q6T3W5
B	10	HIS	-	LINKER	UNP Q6T3W5

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	110	Total	O	0	0
			110	110		

Continued on next page...

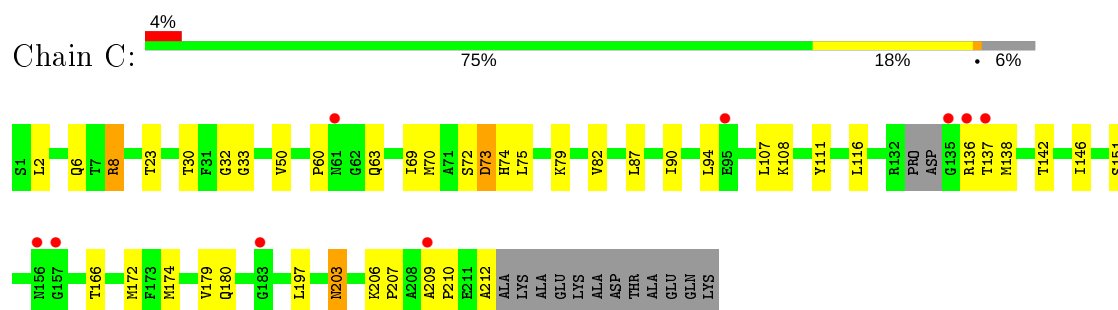
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	106	Total 106	O 106	0	0
4	A	131	Total 131	O 131	0	0
4	B	111	Total 111	O 111	0	0

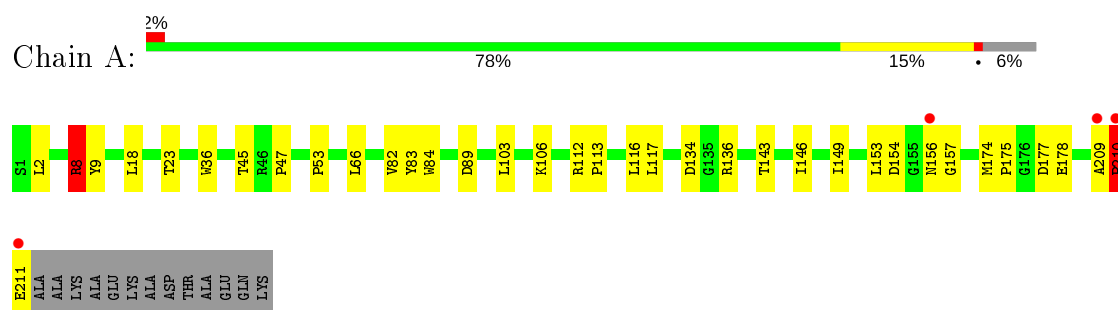
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

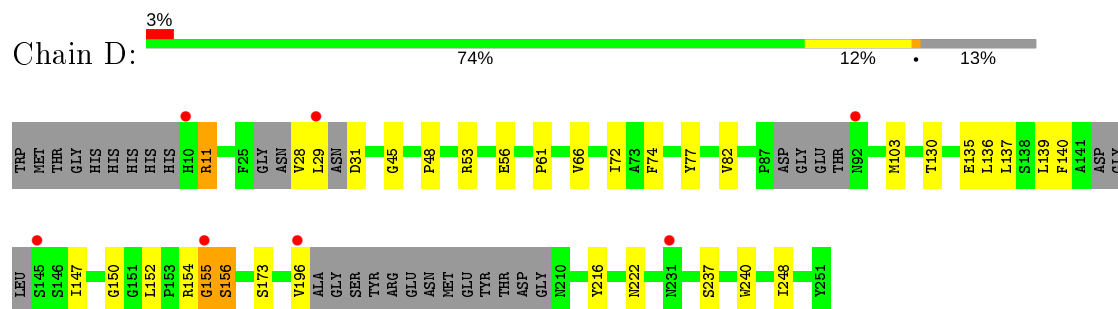
- Molecule 1: Chaperone protein faeE



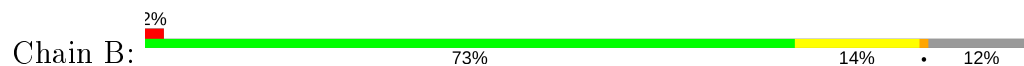
- Molecule 1: Chaperone protein faeE

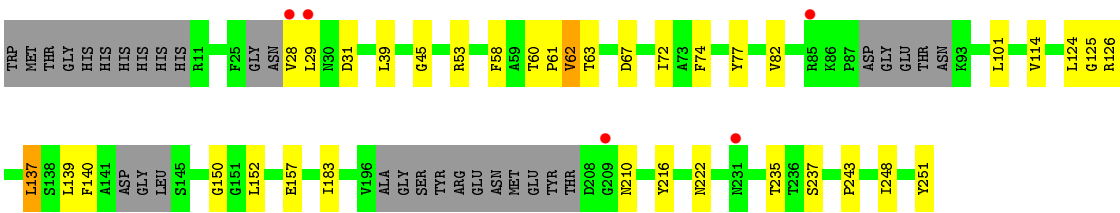


- Molecule 2: FaeG



- Molecule 2: FaeG





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.92Å 90.11Å 86.54Å 90.00° 114.49° 90.00°	Depositor
Resolution (Å)	28.80 – 1.99 28.80 – 1.99	Depositor EDS
% Data completeness (in resolution range)	94.2 (28.80-1.99) 94.4 (28.80-1.99)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 2.00Å)	Xtriage
Refinement program	REFMAC, PHENIX	Depositor
R, R_{free}	0.239 , 0.286 0.242 , 0.282	Depositor DCC
R_{free} test set	3411 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.218	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6775	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.13 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.0841e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.74	1/1623 (0.1%)	0.82	2/2206 (0.1%)
1	C	0.57	0/1587	0.70	1/2157 (0.0%)
2	B	0.52	0/1583	0.67	0/2152
2	D	0.51	0/1574	0.63	0/2136
All	All	0.59	1/6367 (0.0%)	0.71	3/8651 (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	1
2	D	0	2
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	36	TRP	CB-CG	6.09	1.61	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	ARG	CB-CA-C	-6.52	97.36	110.40
1	C	8	ARG	NE-CZ-NH2	-5.78	117.41	120.30
1	A	8	ARG	NE-CZ-NH2	-5.36	117.62	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	PRO	Mainchain
2	D	154	ARG	Peptide
2	D	155	GLY	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	1593	0	1568	27	0
1	C	1559	0	1515	30	0
2	B	1559	0	1512	36	0
2	D	1551	0	1480	25	0
3	A	25	0	0	0	0
3	B	5	0	0	0	0
3	C	20	0	0	0	0
3	D	5	0	0	0	0
4	A	131	0	0	1	0
4	B	111	0	0	1	0
4	C	110	0	0	0	0
4	D	106	0	0	2	0
All	All	6775	0	6075	115	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 (115) 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)
2:D:155:GLY:HA3	2:D:156:SER:CB	1.46	1.41
2:D:155:GLY:CA	2:D:156:SER:CB	2.31	1.06
2:B:60:THR:HG22	2:B:62:VAL:H	1.29	0.97
2:D:45:GLY:H	2:D:222:ASN:HD22	1.09	0.93
2:B:45:GLY:H	2:B:222:ASN:HD22	1.07	0.93
1:C:23:THR:HG22	1:C:63:GLN:HG2	1.62	0.80
1:C:73:ASP:O	1:C:74:HIS:CB	2.30	0.77
2:B:137:LEU:N	2:B:137:LEU:HD23	2.02	0.74
2:D:45:GLY:H	2:D:222:ASN:ND2	1.85	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:45:GLY:N	2:D:222:ASN:HD22	1.85	0.71
1:C:72:SER:O	1:C:73:ASP:O	2.09	0.69
2:B:140:PHE:H	2:B:157:GLU:HG2	1.57	0.68
2:D:72:ILE:HG12	2:D:248:ILE:HG12	1.76	0.68
1:A:209:ALA:O	1:A:211:GLU:N	2.29	0.66
1:C:136:ARG:HB3	1:C:138:MET:HE3	1.78	0.65
1:A:156:ASN:N	1:A:157:GLY:HA2	2.14	0.62
2:D:137:LEU:HD21	2:D:196:VAL:HG23	1.80	0.62
1:A:178:GLU:O	1:A:178:GLU:HG3	2.00	0.62
2:B:126:ARG:HG3	2:B:210:ASN:HD21	1.64	0.62
1:C:73:ASP:OD1	1:C:74:HIS:N	2.28	0.62
2:B:72:ILE:HG12	2:B:248:ILE:HG12	1.83	0.61
1:C:72:SER:C	1:C:73:ASP:O	2.40	0.59
1:C:136:ARG:HB3	1:C:138:MET:CE	2.31	0.59
2:D:103:MET:HE2	2:D:240:TRP:HB2	1.84	0.59
1:C:138:MET:HE2	1:C:180:GLN:HA	1.84	0.58
2:B:124:LEU:HD23	2:B:124:LEU:C	2.25	0.57
2:D:53:ARG:HH21	2:D:173:SER:HA	1.69	0.56
2:D:28:VAL:O	2:D:29:LEU:CB	2.54	0.56
2:B:157:GLU:CD	2:B:157:GLU:H	2.08	0.56
1:A:156:ASN:H	1:A:157:GLY:HA2	1.71	0.55
1:A:53:PRO:HD2	1:A:66:LEU:HD23	1.87	0.55
1:A:82:VAL:C	1:A:83:TYR:CD2	2.80	0.55
2:D:56:GLU:CG	4:D:438:HOH:O	2.54	0.55
2:B:126:ARG:CG	2:B:210:ASN:HD21	2.20	0.54
2:B:63:THR:HG21	2:B:126:ARG:CD	2.37	0.54
1:C:32:GLY:HA3	1:C:90:ILE:HB	1.90	0.54
1:C:209:ALA:HB3	1:C:210:PRO:HD3	1.90	0.54
1:A:154:ASP:OD1	1:A:157:GLY:HA2	2.08	0.53
2:B:28:VAL:O	2:B:31:ASP:HB2	2.09	0.53
2:B:137:LEU:N	2:B:137:LEU:CD2	2.72	0.52
1:A:53:PRO:HD2	1:A:66:LEU:CD2	2.40	0.52
1:A:83:TYR:N	1:A:83:TYR:CD2	2.77	0.52
1:A:134:ASP:OD1	1:A:136:ARG:HB2	2.10	0.52
1:A:9:TYR:HB3	1:A:18:LEU:HD21	1.92	0.52
2:D:136:LEU:O	2:D:137:LEU:HD23	2.09	0.51
1:C:6:GLN:HE22	2:D:11:ARG:HG2	1.76	0.51
2:B:60:THR:HG22	2:B:62:VAL:N	2.11	0.51
2:D:74:PHE:HB2	2:D:82:VAL:O	2.09	0.50
1:A:113:PRO:HD2	1:A:116:LEU:HD12	1.95	0.49
2:B:63:THR:HG21	2:B:126:ARG:HD2	1.94	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:29:LEU:HA	2:D:31:ASP:N	2.27	0.48
2:B:72:ILE:HB	2:B:150:GLY:HA2	1.94	0.48
2:B:29:LEU:HD23	2:B:39:LEU:HD22	1.96	0.48
1:C:108:LYS:HE2	1:C:172:MET:SD	2.53	0.48
1:A:89:ASP:OD2	1:A:103:LEU:HB3	2.13	0.48
2:B:45:GLY:H	2:B:222:ASN:ND2	1.91	0.47
2:B:60:THR:CG2	2:B:62:VAL:O	2.62	0.47
1:A:209:ALA:C	1:A:211:GLU:H	2.18	0.47
2:B:53:ARG:NE	4:B:510:HOH:O	2.47	0.47
2:B:140:PHE:N	2:B:157:GLU:HG2	2.29	0.47
2:D:139:LEU:HA	2:D:140:PHE:HA	1.76	0.47
1:A:45:THR:O	1:A:47:PRO:HD3	2.15	0.47
2:D:135:GLU:HB3	2:D:196:VAL:HG22	1.97	0.47
2:D:77:TYR:HA	4:D:370:HOH:O	2.14	0.47
1:A:112:ARG:HG2	1:A:117:LEU:HD23	1.98	0.46
1:C:70:MET:HE1	1:C:111:TYR:CE1	2.50	0.46
2:B:58:PHE:CE2	2:B:60:THR:OG1	2.69	0.46
1:C:197:LEU:HD23	2:D:61:PRO:HG2	1.98	0.46
1:A:174:MET:HB2	1:A:177:ASP:OD2	2.16	0.46
2:D:103:MET:HE2	2:D:240:TRP:CB	2.46	0.45
2:D:11:ARG:HB2	2:D:11:ARG:NH1	2.31	0.45
2:B:29:LEU:HD23	2:B:29:LEU:HA	1.79	0.45
2:B:74:PHE:HB2	2:B:82:VAL:O	2.16	0.45
2:B:101:LEU:HB2	2:B:114:VAL:HG12	1.98	0.45
2:B:60:THR:HG21	2:B:62:VAL:O	2.17	0.45
1:A:84:TRP:CE2	1:A:106:LYS:HE3	2.52	0.45
1:C:82:VAL:HB	1:C:146:ILE:HD13	1.99	0.44
1:C:70:MET:CE	1:C:111:TYR:CE1	3.01	0.44
1:C:79:LYS:HA	1:C:116:LEU:HD11	1.98	0.44
2:B:139:LEU:HA	2:B:140:PHE:HA	1.72	0.43
2:B:125:GLY:C	2:B:183:ILE:HD11	2.38	0.43
2:D:103:MET:HE3	2:D:240:TRP:CE3	2.52	0.43
1:A:174:MET:HB2	1:A:177:ASP:CG	2.37	0.43
2:B:235:THR:C	2:B:237:SER:H	2.22	0.43
2:B:63:THR:HG21	2:B:126:ARG:HD3	1.99	0.43
1:A:209:ALA:C	1:A:211:GLU:N	2.72	0.43
2:B:60:THR:HG23	2:B:67:ASP:OD2	2.19	0.43
1:A:174:MET:O	1:A:175:PRO:C	2.57	0.43
1:A:2:LEU:HA	1:A:23:THR:O	2.18	0.43
1:C:166:THR:HG23	1:C:179:VAL:HG11	2.01	0.43
2:B:29:LEU:C	2:B:31:ASP:H	2.23	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2:LEU:HB3	1:C:87:LEU:HD21	2.00	0.42
2:D:66:VAL:N	2:D:147:ILE:HD12	2.34	0.42
2:B:77:TYR:HB3	2:B:243:PRO:HG2	2.00	0.42
1:C:50:VAL:HG22	1:C:69:ILE:CD1	2.49	0.42
1:C:137:THR:O	1:C:138:MET:HE2	2.20	0.42
1:C:82:VAL:HG11	1:C:174:MET:HE1	2.01	0.42
1:A:143:THR:O	1:A:211:GLU:OE1	2.38	0.42
1:C:172:MET:HE2	1:C:172:MET:HB3	1.87	0.42
2:B:58:PHE:HE2	2:B:60:THR:OG1	2.03	0.41
1:C:30:THR:OG1	1:C:60:PRO:HD3	2.21	0.41
1:C:94:LEU:HD23	1:C:94:LEU:HA	1.66	0.41
2:D:137:LEU:HD21	2:D:196:VAL:CG2	2.49	0.41
4:A:353:HOH:O	2:B:61:PRO:HB2	2.20	0.41
1:C:50:VAL:HG22	1:C:69:ILE:HG13	2.02	0.41
1:A:84:TRP:CD2	1:A:106:LYS:HE3	2.55	0.41
1:A:8:ARG:NH2	2:B:251:TYR:OXT	2.46	0.41
1:C:2:LEU:HD11	1:C:33:GLY:HA3	2.02	0.41
2:D:72:ILE:HB	2:D:150:GLY:HA2	2.02	0.41
1:C:142:THR:HG22	1:C:212:ALA:HA	2.03	0.41
1:C:166:THR:HG23	1:C:179:VAL:CG1	2.50	0.41
1:A:82:VAL:HB	1:A:146:ILE:HD13	2.03	0.40
1:C:206:LYS:HA	1:C:207:PRO:HD2	1.95	0.40
1:A:209:ALA:O	1:A:210:PRO:C	2.57	0.40
2:B:29:LEU:C	2:B:31:ASP:N	2.75	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	210/224 (94%)	201 (96%)	8 (4%)	1 (0%)	29 23

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	206/224 (92%)	199 (97%)	5 (2%)	2 (1%)	15	9
2	B	210/251 (84%)	205 (98%)	5 (2%)	0	100	100
2	D	208/251 (83%)	202 (97%)	5 (2%)	1 (0%)	29	23
All	All	834/950 (88%)	807 (97%)	23 (3%)	4 (0%)	29	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	73	ASP
1	C	203	ASN
2	D	156	SER
1	A	210	PRO

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	164/189 (87%)	161 (98%)	3 (2%)	59	63
1	C	157/189 (83%)	152 (97%)	5 (3%)	39	38
2	B	152/192 (79%)	148 (97%)	4 (3%)	46	48
2	D	148/192 (77%)	142 (96%)	6 (4%)	30	28
All	All	621/762 (82%)	603 (97%)	18 (3%)	42	43

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	8	ARG
1	C	75	LEU
1	C	107	LEU
1	C	151	SER
1	C	203	ASN
2	D	11	ARG
2	D	48	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	D	130	THR
2	D	152	LEU
2	D	216	TYR
2	D	237	SER
1	A	8	ARG
1	A	149	ILE
1	A	153	LEU
2	B	62	VAL
2	B	137	LEU
2	B	152	LEU
2	B	216	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	6	GLN
1	C	203	ASN
2	D	222	ASN
2	B	210	ASN
2	B	222	ASN

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](#)

11 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	$\# Z > 2$
3	SO4	C	228	-	4,4,4	0.16	0	6,6,6	0.21	0
3	SO4	A	227	-	4,4,4	0.10	0	6,6,6	0.23	0
3	SO4	C	227	-	4,4,4	0.17	0	6,6,6	0.24	0
3	SO4	C	226	-	4,4,4	0.10	0	6,6,6	0.18	0
3	SO4	A	228	-	4,4,4	0.13	0	6,6,6	0.40	0
3	SO4	A	226	-	4,4,4	0.14	0	6,6,6	0.17	0
3	SO4	B	252	-	4,4,4	0.21	0	6,6,6	0.27	0
3	SO4	D	252	-	4,4,4	0.24	0	6,6,6	0.16	0
3	SO4	C	225	-	4,4,4	0.18	0	6,6,6	0.33	0
3	SO4	A	225	-	4,4,4	0.18	0	6,6,6	0.32	0
3	SO4	A	229	-	4,4,4	0.21	0	6,6,6	0.26	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 ⓘ

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	211/224 (94%)	-0.02	4 (1%) 66 65	13, 24, 42, 61	0
1	C	210/224 (93%)	0.12	9 (4%) 35 34	16, 29, 54, 65	0
2	B	220/251 (87%)	0.01	5 (2%) 60 59	15, 27, 44, 64	0
2	D	219/251 (87%)	0.10	7 (3%) 47 46	16, 28, 49, 68	0
All	All	860/950 (90%)	0.05	25 (2%) 51 50	13, 27, 49, 68	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	29	LEU	6.3
2	D	10	HIS	5.4
2	B	28	VAL	4.5
2	D	155	GLY	4.3
1	C	209	ALA	4.1
1	A	211	GLU	4.1
1	C	156	ASN	3.8
1	A	156	ASN	3.4
1	A	209	ALA	3.3
2	B	209	GLY	3.1
2	D	29	LEU	2.7
1	A	210	PRO	2.6
2	B	85	ARG	2.6
2	D	145	SER	2.5
1	C	136	ARG	2.5
1	C	137	THR	2.5
2	D	92	ASN	2.5
2	D	231	ASN	2.5
1	C	183	GLY	2.4
2	D	196	VAL	2.4
1	C	135	GLY	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	231	ASN	2.2
1	C	95	GLU	2.2
1	C	61	ASN	2.1
1	C	157	GLY	2.1

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. 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(\AA^2)	Q<0.9
3	SO4	A	228	5/5	0.81	0.19	74,77,77,79	0
3	SO4	C	228	5/5	0.84	0.17	75,76,77,78	0
3	SO4	A	226	5/5	0.87	0.10	88,88,89,89	0
3	SO4	A	225	5/5	0.91	0.19	66,66,67,70	0
3	SO4	C	225	5/5	0.92	0.20	55,56,58,60	0
3	SO4	A	227	5/5	0.96	0.08	50,51,53,54	0
3	SO4	A	229	5/5	0.96	0.16	42,45,48,50	0
3	SO4	C	226	5/5	0.97	0.09	51,52,53,54	0
3	SO4	C	227	5/5	0.97	0.16	56,57,57,58	0
3	SO4	D	252	5/5	0.98	0.08	36,38,41,42	0
3	SO4	B	252	5/5	0.98	0.08	37,40,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.