



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 05:52 PM BST

PDB ID : 6SDB
Title : Chimeric titin Z1Z2 functionalized with a KLER exogenous peptide from decorin
Authors : Fleming, J.R.; Hill, C.; Mayans, O.M.
Deposited on : 2019-07-26
Resolution : 2.80 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 : 2.13.1
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) : 2.13.1

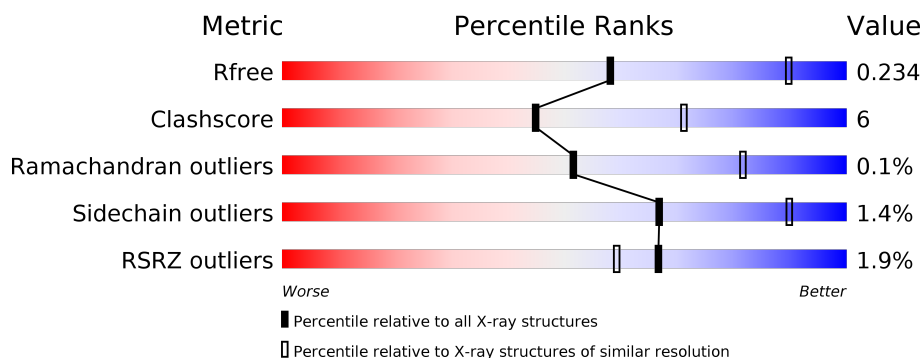
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 85%, yellow 14%, grey 0%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 85% 14% . </div> </div>
1	B	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 81%, yellow 18%, grey 0%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 81% 18% .. </div> </div>
1	C	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, green 42%, yellow 8%, grey 50%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 42% 8% . 50% </div> </div>
1	D	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, green 41%, yellow 8%, grey 51%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 41% 8% 51% </div> </div>
1	E	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 85%, yellow 13%, grey 0%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 85% 13% .. </div> </div>

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
4	EDO	B	203	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6141 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 Titin,Decorin,Titin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1470	921	254	294	1			
1	B	196	Total	C	N	O	S	0	0	0
			1473	923	255	293	2			
1	C	99	Total	C	N	O	S	1	0	0
			738	466	127	144	1			
1	D	97	Total	C	N	O	S	2	0	0
			740	460	128	151	1			
1	E	193	Total	C	N	O	S	0	0	0
			1460	916	252	290	2			

There are 20 discrepancies between the modelled and reference sequences:

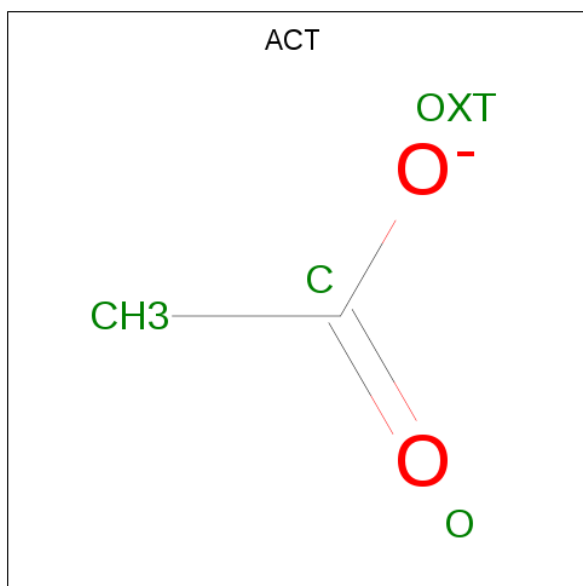
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q8WZ42
A	0	ALA	-	expression tag	UNP Q8WZ42
A	1	MET	-	expression tag	UNP Q8WZ42
A	2	ALA	-	expression tag	UNP Q8WZ42
B	-1	GLY	-	expression tag	UNP Q8WZ42
B	0	ALA	-	expression tag	UNP Q8WZ42
B	1	MET	-	expression tag	UNP Q8WZ42
B	2	ALA	-	expression tag	UNP Q8WZ42
C	-1	GLY	-	expression tag	UNP Q8WZ42
C	0	ALA	-	expression tag	UNP Q8WZ42
C	1	MET	-	expression tag	UNP Q8WZ42
C	2	ALA	-	expression tag	UNP Q8WZ42
D	-1	GLY	-	expression tag	UNP Q8WZ42
D	0	ALA	-	expression tag	UNP Q8WZ42
D	1	MET	-	expression tag	UNP Q8WZ42
D	2	ALA	-	expression tag	UNP Q8WZ42
E	-1	GLY	-	expression tag	UNP Q8WZ42
E	0	ALA	-	expression tag	UNP Q8WZ42
E	1	MET	-	expression tag	UNP Q8WZ42

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Chain	Residue	Modelled	Actual	Comment	Reference
E	2	ALA	-	expression tag	UNP Q8WZ42

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



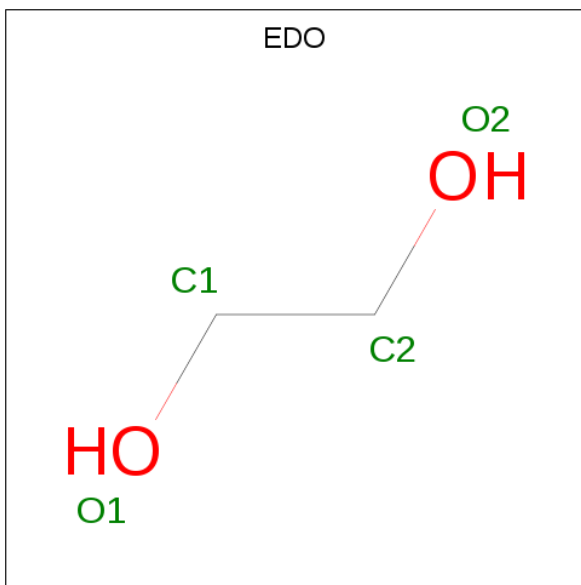
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total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

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



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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is water.

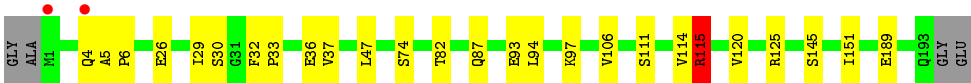
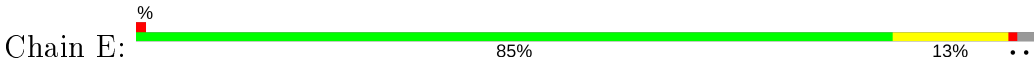
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	44	Total	O	0	0
			44	44		
6	B	53	Total	O	0	0
			53	53		
6	C	24	Total	O	0	0
			24	24		
6	D	17	Total	O	0	0
			17	17		
6	E	64	Total	O	0	0
			64	64		

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.

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|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| GLY | ALA | MET | ALA | ALA | THR | GLN | ALA | ALA | PRO | THR | PHE | THR | GLN | PRO | LEU | GLY | SER | VAL | VAL | VAL | VAL | LEU | LEU | GLU | GLY | SER | THR | ALA | ALA | PHE | THR | THR | GLU | ALA | HIS | ILE | ILE | SER | SER | GLY | PHE | PHE | PRO | PRO | VAL | VAL | VAL | PRO | GLU | GLU | VAL | VAL | SER | SER | TRP | PHE | ARG | ASP | GLY | GLN | GLN | VAL | LYS | LEU | LEU | GLU | GLU | GLN | GLN | ILE | ILE | SER | SER | PHE | PHE |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|



● Molecule 1: Titin,Decorin,Titin



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	149.20Å 149.20Å 149.91Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.84 – 2.80 48.94 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.84-2.80) 99.9 (48.94-2.80)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.13_2998, PHENIX 1.13_2998	Depositor
R, R_{free}	0.191 , 0.234 0.190 , 0.234	Depositor DCC
R_{free} test set	1781 reflections (3.82%)	wwPDB-VP
Wilson B-factor (Å ²)	64.2	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.035 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6141	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% 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, 0.333 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: PO4, ACT, EDO, 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	0.51	0/1496	0.71	1/2037 (0.0%)
1	B	0.52	0/1499	0.88	5/2040 (0.2%)
1	C	0.53	1/752 (0.1%)	0.72	1/1022 (0.1%)
1	D	0.42	0/751	0.63	0/1022
1	E	0.49	0/1486	0.73	4/2023 (0.2%)
All	All	0.50	1/5984 (0.0%)	0.75	11/8144 (0.1%)

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	E	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	93	GLU	CB-CG	-5.51	1.41	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	115	ARG	NE-CZ-NH1	-11.81	114.39	120.30
1	B	115	ARG	NE-CZ-NH2	9.55	125.08	120.30
1	B	115	ARG	CG-CD-NE	7.94	128.48	111.80
1	A	115	ARG	NE-CZ-NH2	-6.96	116.82	120.30
1	B	61	ARG	CB-CG-CD	-6.73	94.10	111.60
1	E	26	GLU	N-CA-CB	6.16	121.68	110.60
1	E	36	GLU	N-CA-CB	-5.74	100.26	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	115	ARG	CD-NE-CZ	5.54	131.36	123.60
1	B	115	ARG	CD-NE-CZ	-5.54	115.85	123.60
1	C	76	ARG	CG-CD-NE	-5.47	100.31	111.80
1	E	26	GLU	CB-CA-C	-5.42	99.56	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	115	ARG	Sidechain

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	1470	0	1463	19	0
1	B	1473	0	1472	30	0
1	C	738	0	746	8	0
1	D	740	0	725	7	0
1	E	1460	0	1461	14	0
2	A	24	0	18	1	0
2	B	4	0	3	0	0
2	D	4	0	3	0	0
2	E	12	0	9	0	0
3	B	5	0	0	1	0
4	B	4	0	6	0	0
5	E	5	0	0	0	0
6	A	44	0	0	0	0
6	B	53	0	0	1	0
6	C	24	0	0	0	0
6	D	17	0	0	0	0
6	E	64	0	0	1	0
All	All	6141	0	5906	77	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 6.

All (77) 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:B:115:ARG:HG2	1:B:193:GLN:HG3	1.43	0.98
1:D:138:ARG:NH1	1:D:170:SER:O	2.18	0.77
1:E:111:SER:HB3	1:E:189:GLU:H	1.49	0.76
1:B:76:ARG:CZ	1:B:91:THR:HG21	2.17	0.74
1:B:37:VAL:HB	1:B:79:LEU:HD11	1.72	0.71
1:A:114:VAL:HG22	1:A:192:VAL:HG12	1.71	0.71
1:A:47:LEU:HD12	1:A:50:LEU:HD12	1.73	0.70
1:A:93:GLU:HB2	2:A:204:ACT:H1	1.77	0.65
1:A:82:THR:HG23	1:A:87:GLN:HB2	1.79	0.64
1:B:120:VAL:HG22	1:B:162:ILE:HB	1.81	0.63
1:D:115:ARG:HD3	1:D:116:GLN:H	1.65	0.62
1:E:82:THR:HG22	1:E:87:GLN:HG3	1.81	0.61
1:B:98:ALA:N	3:B:202:SO4:O1	2.27	0.61
1:E:4:GLN:HG3	1:E:5:ALA:N	2.17	0.60
1:B:47:LEU:HD12	1:B:50:LEU:HD12	1.84	0.59
1:B:167:PRO:HA	1:B:192:VAL:HG22	1.86	0.58
1:B:107:GLN:HB3	1:B:125:ARG:HG3	1.86	0.58
1:B:36:GLU:HG2	1:B:37:VAL:N	2.20	0.56
1:B:16:VAL:HG12	1:B:129:ILE:HD11	1.88	0.55
1:B:97:LYS:NZ	6:B:301:HOH:O	2.37	0.55
1:B:36:GLU:O	1:B:81:ALA:HA	2.07	0.55
1:A:25:PHE:HZ	1:A:94:LEU:HB2	1.74	0.53
1:A:166:TYR:O	1:A:192:VAL:HG21	2.09	0.52
1:B:114:VAL:HG22	1:B:192:VAL:HG12	1.92	0.52
1:B:16:VAL:HG22	1:B:95:LEU:HB3	1.92	0.51
1:A:120:VAL:HG13	1:A:162:ILE:HB	1.92	0.51
1:A:167:PRO:HA	1:A:192:VAL:HG23	1.93	0.51
1:E:114:VAL:HG11	1:E:120:VAL:HG22	1.93	0.51
1:E:93:GLU:OE2	1:E:93:GLU:N	2.44	0.50
1:C:76:ARG:HH21	1:C:91:THR:HG21	1.77	0.49
1:B:71:LYS:HE2	1:B:98:ALA:HB2	1.93	0.49
1:D:151:ILE:HA	1:D:159:SER:O	2.13	0.49
1:A:114:VAL:O	1:A:192:VAL:HA	2.12	0.49
1:C:76:ARG:HE	1:C:91:THR:HG21	1.76	0.49
1:E:4:GLN:NE2	6:E:301:HOH:O	2.45	0.49
1:E:6:PRO:HA	1:E:30:SER:O	2.12	0.49
1:C:46:LYS:HB2	1:C:49:ARG:HG3	1.94	0.48
1:E:106:VAL:HG22	1:E:125:ARG:HB3	1.95	0.48
1:B:76:ARG:HG3	1:B:93:GLU:OE2	2.13	0.48
1:D:167:PRO:HA	1:D:192:VAL:HB	1.96	0.47
1:D:114:VAL:HG11	1:D:120:VAL:HG22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:THR:HG22	1:B:129:ILE:HG22	1.97	0.47
1:B:36:GLU:HG2	1:B:37:VAL:H	1.79	0.47
1:E:145:SER:HB2	1:E:151:ILE:HB	1.97	0.47
1:A:37:VAL:HA	1:A:80:LYS:O	2.15	0.46
1:B:18:LEU:HB2	1:B:129:ILE:HG21	1.96	0.46
1:B:47:LEU:HD12	1:B:47:LEU:HA	1.79	0.46
1:C:47:LEU:HD12	1:C:50:LEU:HD12	1.97	0.46
1:A:167:PRO:HA	1:A:192:VAL:CG2	2.46	0.46
1:A:165:ALA:HB1	1:A:192:VAL:HG11	1.98	0.46
1:A:54:GLN:HG2	1:B:108:ARG:NH1	2.31	0.46
1:B:115:ARG:HD3	1:B:115:ARG:HH11	1.32	0.46
1:B:32:PHE:HA	1:B:33:PRO:C	2.37	0.45
1:B:166:TYR:C	1:B:192:VAL:HG21	2.36	0.45
1:A:114:VAL:CG2	1:A:192:VAL:HG12	2.41	0.45
1:D:103:PRO:HA	1:D:127:THR:O	2.16	0.45
1:E:47:LEU:HA	1:E:47:LEU:HD23	1.75	0.45
1:A:12:LEU:HD22	1:A:25:PHE:HB3	1.99	0.45
1:E:106:VAL:CG2	1:E:125:ARG:HB3	2.47	0.44
1:E:32:PHE:HA	1:E:33:PRO:C	2.38	0.44
1:B:144:GLN:H	1:B:144:GLN:CD	2.21	0.44
1:E:29:ILE:HD12	1:E:37:VAL:HG11	2.00	0.44
1:A:138:ARG:NH2	1:A:170:SER:O	2.51	0.43
1:A:113:THR:HA	1:A:191:LEU:O	2.18	0.43
1:D:145:SER:HB2	1:D:151:ILE:HG12	2.01	0.43
1:C:7:THR:HG22	1:C:30:SER:OG	2.18	0.42
1:A:25:PHE:CZ	1:A:94:LEU:HB2	2.55	0.42
1:B:108:ARG:HH11	1:B:108:ARG:HG2	1.84	0.42
1:B:95:LEU:HD21	1:B:97:LYS:HE3	2.01	0.42
1:E:74:SER:HA	1:E:94:LEU:O	2.20	0.42
1:B:167:PRO:HA	1:B:192:VAL:CG2	2.49	0.41
1:C:37:VAL:HA	1:C:80:LYS:O	2.20	0.41
1:B:7:THR:HG23	1:B:30:SER:HB2	2.01	0.41
1:C:82:THR:OG1	1:C:87:GLN:HG3	2.21	0.41
1:A:32:PHE:HA	1:A:33:PRO:C	2.40	0.41
1:C:32:PHE:HA	1:C:33:PRO:C	2.41	0.40
1:B:113:THR:HA	1:B:191:LEU:O	2.20	0.40

There are no symmetry-related clashes.

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	193/197 (98%)	192 (100%)	0	1 (0%)	29	61
1	B	194/197 (98%)	192 (99%)	2 (1%)	0	100	100
1	C	97/197 (49%)	97 (100%)	0	0	100	100
1	D	95/197 (48%)	94 (99%)	1 (1%)	0	100	100
1	E	191/197 (97%)	190 (100%)	1 (0%)	0	100	100
All	All	770/985 (78%)	765 (99%)	4 (0%)	1 (0%)	51	81

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	GLY

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	161/162 (99%)	159 (99%)	2 (1%)	71	92
1	B	161/162 (99%)	160 (99%)	1 (1%)	86	96
1	C	80/162 (49%)	79 (99%)	1 (1%)	69	91
1	D	82/162 (51%)	79 (96%)	3 (4%)	34	68
1	E	161/162 (99%)	159 (99%)	2 (1%)	71	92
All	All	645/810 (80%)	636 (99%)	9 (1%)	67	90

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	ARG
1	A	95	LEU
1	B	119	GLN
1	C	90	SER
1	D	108	ARG
1	D	125	ARG
1	D	157	LEU
1	E	97	LYS
1	E	115	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	116	GLN
1	B	123	GLN
1	C	13	GLN

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](#)

14 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	ACT	A	204	-	1,3,3	9.03	1 (100%)	0,3,3	0.00	-
5	PO4	E	201	-	4,4,4	0.85	0	6,6,6	0.54	0
2	ACT	A	203	-	1,3,3	9.83	1 (100%)	0,3,3	0.00	-
4	EDO	B	203	-	3,3,3	0.64	0	2,2,2	0.30	0
2	ACT	E	203	-	1,3,3	9.60	1 (100%)	0,3,3	0.00	-
2	ACT	E	204	-	1,3,3	9.74	1 (100%)	0,3,3	0.00	-
2	ACT	A	202	-	1,3,3	8.33	1 (100%)	0,3,3	0.00	-
3	SO4	B	202	-	4,4,4	0.14	0	6,6,6	0.19	0
2	ACT	E	202	-	1,3,3	8.76	1 (100%)	0,3,3	0.00	-
2	ACT	A	201	-	1,3,3	5.60	1 (100%)	0,3,3	0.00	-
2	ACT	B	201	-	1,3,3	8.87	1 (100%)	0,3,3	0.00	-
2	ACT	D	201	-	1,3,3	9.38	1 (100%)	0,3,3	0.00	-
2	ACT	A	205	-	1,3,3	9.42	1 (100%)	0,3,3	0.00	-
2	ACT	A	206	-	1,3,3	9.84	1 (100%)	0,3,3	0.00	-

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	B	203	-	-	0/1/1/1	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	206	ACT	CH3-C	9.84	1.61	1.48
2	A	203	ACT	CH3-C	9.83	1.61	1.48
2	E	204	ACT	CH3-C	9.74	1.61	1.48
2	E	203	ACT	CH3-C	9.60	1.61	1.48
2	A	205	ACT	CH3-C	9.42	1.60	1.48
2	D	201	ACT	CH3-C	9.38	1.60	1.48
2	A	204	ACT	CH3-C	9.03	1.60	1.48
2	B	201	ACT	CH3-C	8.87	1.60	1.48
2	E	202	ACT	CH3-C	8.76	1.59	1.48
2	A	202	ACT	CH3-C	8.33	1.59	1.48
2	A	201	ACT	CH3-C	5.60	1.55	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	204	ACT	1	0
3	B	202	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	195/197 (98%)	-0.27	2 (1%) 82 77	43, 58, 86, 129	0
1	B	196/197 (99%)	-0.23	1 (0%) 91 88	44, 60, 100, 149	0
1	C	99/197 (50%)	-0.02	4 (4%) 38 28	49, 62, 97, 129	1 (1%)
1	D	97/197 (49%)	0.11	6 (6%) 20 13	50, 75, 105, 130	1 (1%)
1	E	193/197 (97%)	-0.17	2 (1%) 82 77	44, 62, 96, 125	0
All	All	780/985 (79%)	-0.15	15 (1%) 66 59	43, 62, 101, 149	2 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1	MET	4.1
1	D	101	ALA	3.6
1	C	99	GLU	3.4
1	B	147	LEU	3.3
1	A	1	MET	3.0
1	D	156	ASP	2.9
1	C	2	ALA	2.5
1	D	132	PRO	2.4
1	D	107	GLN	2.4
1	D	100	THR	2.4
1	D	155	GLY	2.3
1	C	18	LEU	2.3
1	A	2	ALA	2.3
1	E	4	GLN	2.3
1	E	1	MET	2.2

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
4	EDO	B	203	4/4	0.74	0.45	73,86,87,90	0
2	ACT	D	201	4/4	0.79	0.37	90,90,96,99	0
2	ACT	B	201	4/4	0.82	0.33	91,92,94,95	0
2	ACT	E	202	4/4	0.88	0.58	77,77,78,80	0
2	ACT	A	205	4/4	0.90	0.50	79,91,92,98	0
2	ACT	A	202	4/4	0.91	0.60	84,93,94,97	0
2	ACT	A	204	4/4	0.92	0.26	91,92,96,99	0
2	ACT	E	204	4/4	0.93	0.37	46,62,63,64	0
5	PO4	E	201	5/5	0.93	0.17	120,123,129,130	0
2	ACT	A	206	4/4	0.93	0.36	45,56,66,73	0
2	ACT	A	203	4/4	0.94	0.44	58,72,74,80	0
3	SO4	B	202	5/5	0.96	0.19	115,119,121,127	0
2	ACT	A	201	4/4	0.97	0.14	44,52,53,61	0
2	ACT	E	203	4/4	0.97	0.15	61,65,68,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.