



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 02:19 am GMT

PDB ID : 3EZQ
Title : Crystal Structure of the Fas/FADD Death Domain Complex
Authors : Schwarzenbacher, R.; Robinson, H.; Stec, B.; Riedl, S.J.
Deposited on : 2008-10-23
Resolution : 2.73 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

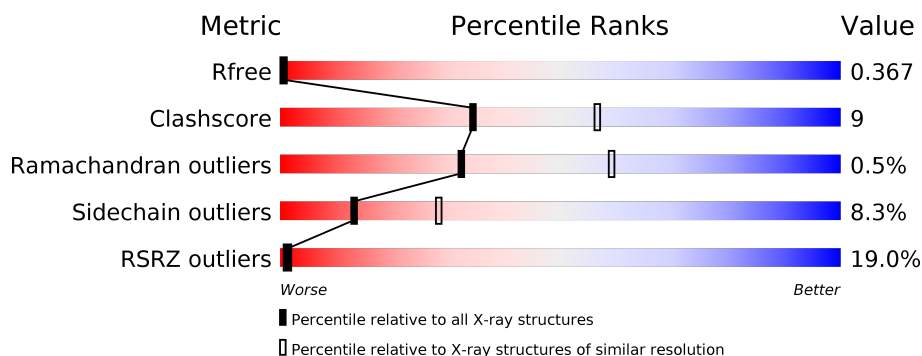
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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}	100719	3342 (2.78-2.70)
Clashscore	112137	3731 (2.78-2.70)
Ramachandran outliers	110173	3670 (2.78-2.70)
Sidechain outliers	110143	3671 (2.78-2.70)
RSRZ outliers	101464	3362 (2.78-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	115	<div> <div>13%</div> <div>81%</div> <div>18%</div> <div>.</div> </div>
1	C	115	<div> <div>17%</div> <div>80%</div> <div>16%</div> <div>.</div> </div>
1	E	115	<div> <div>13%</div> <div>82%</div> <div>17%</div> <div>.</div> </div>
1	G	115	<div> <div>17%</div> <div>87%</div> <div>11%</div> <div>.</div> </div>
1	I	115	<div> <div>16%</div> <div>87%</div> <div>10%</div> <div>.</div> </div>
1	K	115	<div> <div>13%</div> <div>77%</div> <div>20%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
1	M	115	
1	O	115	
2	B	122	
2	D	122	
2	F	122	
2	H	122	
2	J	122	
2	L	122	
2	N	122	
2	P	122	

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
3	SO4	K	13	-	-	X	-
4	NA	B	21	-	-	-	X
4	NA	E	14	-	-	-	X
4	NA	K	16	-	-	-	X
4	NA	M	17	-	-	-	X
4	NA	P	18	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13933 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 Tumor necrosis factor receptor superfamily member 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			
1	C	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			
1	E	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			
1	G	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			
1	I	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			
1	K	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			
1	M	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			
1	O	115	Total	C	N	O	S	0	0	0
			916	571	159	184	2			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	336	LEU	-	EXPRESSION TAG	UNP P25445
A	337	GLU	-	EXPRESSION TAG	UNP P25445
C	336	LEU	-	EXPRESSION TAG	UNP P25445
C	337	GLU	-	EXPRESSION TAG	UNP P25445
E	336	LEU	-	EXPRESSION TAG	UNP P25445
E	337	GLU	-	EXPRESSION TAG	UNP P25445
G	336	LEU	-	EXPRESSION TAG	UNP P25445
G	337	GLU	-	EXPRESSION TAG	UNP P25445
I	336	LEU	-	EXPRESSION TAG	UNP P25445
I	337	GLU	-	EXPRESSION TAG	UNP P25445
K	336	LEU	-	EXPRESSION TAG	UNP P25445
K	337	GLU	-	EXPRESSION TAG	UNP P25445
M	336	LEU	-	EXPRESSION TAG	UNP P25445

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Chain	Residue	Modelled	Actual	Comment	Reference
M	337	GLU	-	EXPRESSION TAG	UNP P25445
O	336	LEU	-	EXPRESSION TAG	UNP P25445
O	337	GLU	-	EXPRESSION TAG	UNP P25445

- Molecule 2 is a protein called Protein FADD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			
2	D	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			
2	F	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			
2	H	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			
2	J	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			
2	L	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			
2	N	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			
2	P	99	Total	C	N	O	S	0	0	0
			795	483	154	154	4			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	209	HIS	-	EXPRESSION TAG	UNP Q13158
B	210	HIS	-	EXPRESSION TAG	UNP Q13158
B	211	HIS	-	EXPRESSION TAG	UNP Q13158
B	212	HIS	-	EXPRESSION TAG	UNP Q13158
B	213	HIS	-	EXPRESSION TAG	UNP Q13158
B	214	HIS	-	EXPRESSION TAG	UNP Q13158
D	209	HIS	-	EXPRESSION TAG	UNP Q13158
D	210	HIS	-	EXPRESSION TAG	UNP Q13158
D	211	HIS	-	EXPRESSION TAG	UNP Q13158
D	212	HIS	-	EXPRESSION TAG	UNP Q13158
D	213	HIS	-	EXPRESSION TAG	UNP Q13158
D	214	HIS	-	EXPRESSION TAG	UNP Q13158
F	209	HIS	-	EXPRESSION TAG	UNP Q13158
F	210	HIS	-	EXPRESSION TAG	UNP Q13158
F	211	HIS	-	EXPRESSION TAG	UNP Q13158

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Chain	Residue	Modelled	Actual	Comment	Reference
F	212	HIS	-	EXPRESSION TAG	UNP Q13158
F	213	HIS	-	EXPRESSION TAG	UNP Q13158
F	214	HIS	-	EXPRESSION TAG	UNP Q13158
H	209	HIS	-	EXPRESSION TAG	UNP Q13158
H	210	HIS	-	EXPRESSION TAG	UNP Q13158
H	211	HIS	-	EXPRESSION TAG	UNP Q13158
H	212	HIS	-	EXPRESSION TAG	UNP Q13158
H	213	HIS	-	EXPRESSION TAG	UNP Q13158
H	214	HIS	-	EXPRESSION TAG	UNP Q13158
J	209	HIS	-	EXPRESSION TAG	UNP Q13158
J	210	HIS	-	EXPRESSION TAG	UNP Q13158
J	211	HIS	-	EXPRESSION TAG	UNP Q13158
J	212	HIS	-	EXPRESSION TAG	UNP Q13158
J	213	HIS	-	EXPRESSION TAG	UNP Q13158
J	214	HIS	-	EXPRESSION TAG	UNP Q13158
L	209	HIS	-	EXPRESSION TAG	UNP Q13158
L	210	HIS	-	EXPRESSION TAG	UNP Q13158
L	211	HIS	-	EXPRESSION TAG	UNP Q13158
L	212	HIS	-	EXPRESSION TAG	UNP Q13158
L	213	HIS	-	EXPRESSION TAG	UNP Q13158
L	214	HIS	-	EXPRESSION TAG	UNP Q13158
N	209	HIS	-	EXPRESSION TAG	UNP Q13158
N	210	HIS	-	EXPRESSION TAG	UNP Q13158
N	211	HIS	-	EXPRESSION TAG	UNP Q13158
N	212	HIS	-	EXPRESSION TAG	UNP Q13158
N	213	HIS	-	EXPRESSION TAG	UNP Q13158
N	214	HIS	-	EXPRESSION TAG	UNP Q13158
P	209	HIS	-	EXPRESSION TAG	UNP Q13158
P	210	HIS	-	EXPRESSION TAG	UNP Q13158
P	211	HIS	-	EXPRESSION TAG	UNP Q13158
P	212	HIS	-	EXPRESSION TAG	UNP Q13158
P	213	HIS	-	EXPRESSION TAG	UNP Q13158
P	214	HIS	-	EXPRESSION TAG	UNP Q13158

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	J	1	Total	O	S	0	0
			5	4	1		
3	K	1	Total	O	S	0	0
			5	4	1		
3	K	1	Total	O	S	0	0
			5	4	1		
3	M	1	Total	O	S	0	0
			5	4	1		
3	M	1	Total	O	S	0	0
			5	4	1		
3	O	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	P	1	Total Na 1 1	0	0
4	J	1	Total Na 1 1	0	0
4	K	1	Total Na 1 1	0	0
4	E	1	Total Na 1 1	0	0
4	H	1	Total Na 1 1	0	0
4	B	1	Total Na 1 1	0	0
4	A	1	Total Na 1 1	0	0
4	M	1	Total Na 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	7	Total O 7 7	0	0
5	B	2	Total O 2 2	0	0
5	C	8	Total O 8 8	0	0
5	D	4	Total O 4 4	0	0
5	E	9	Total O 9 9	0	0
5	F	4	Total O 4 4	0	0
5	G	11	Total O 11 11	0	0
5	H	8	Total O 8 8	0	0
5	I	84	Total O 84 84	0	0
5	J	8	Total O 8 8	0	0
5	K	6	Total O 6 6	0	0
5	L	7	Total O 7 7	0	0

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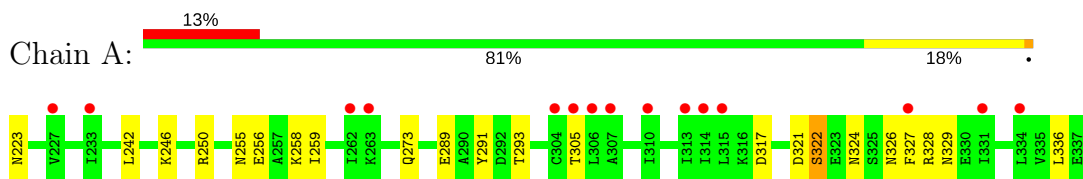
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	M	9	Total 9	O 9	0	0
5	N	3	Total 3	O 3	0	0
5	O	5	Total 5	O 5	0	0
5	P	2	Total 2	O 2	0	0

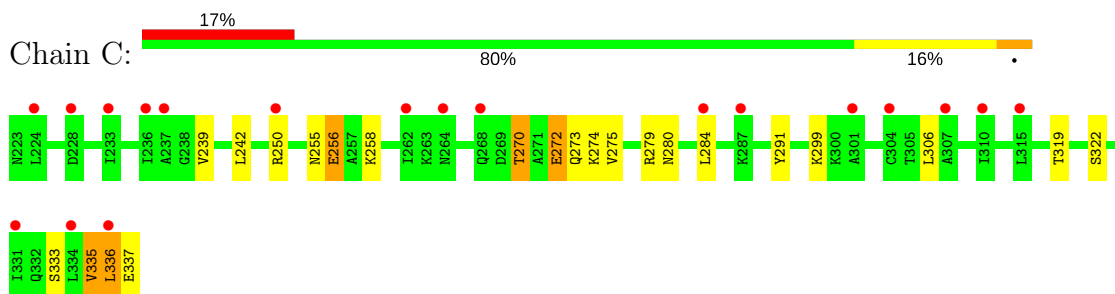
3 Residue-property plots [i](#)

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.

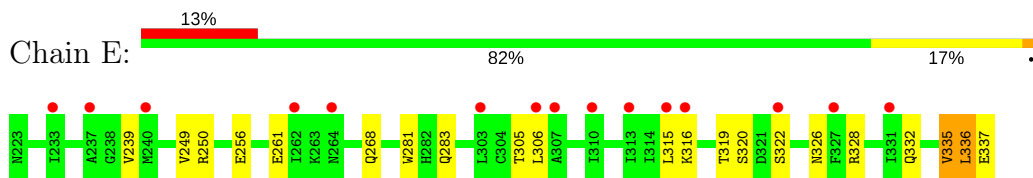
- Molecule 1: Tumor necrosis factor receptor superfamily member 6



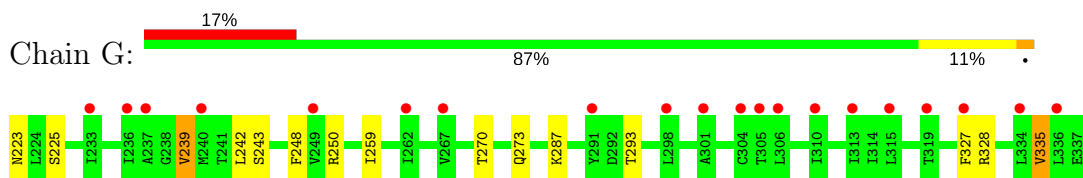
- Molecule 1: Tumor necrosis factor receptor superfamily member 6



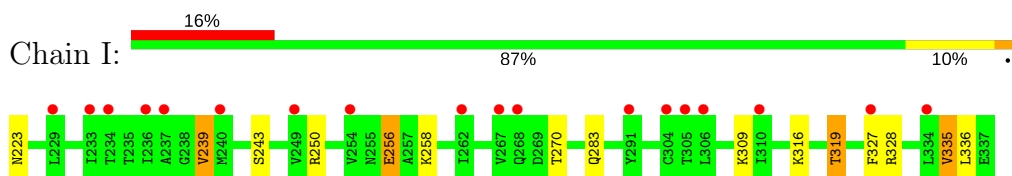
- Molecule 1: Tumor necrosis factor receptor superfamily member 6

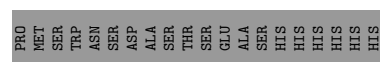
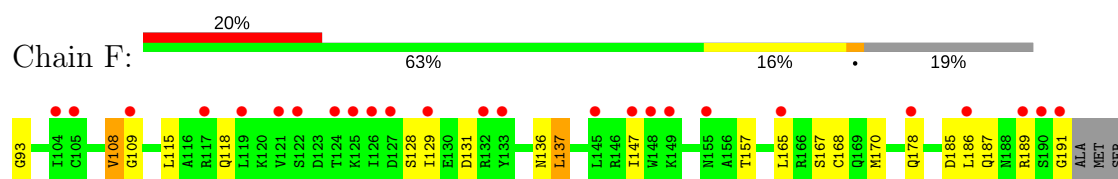


- Molecule 1: Tumor necrosis factor receptor superfamily member 6

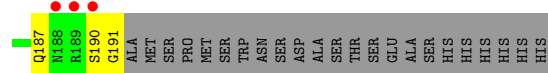
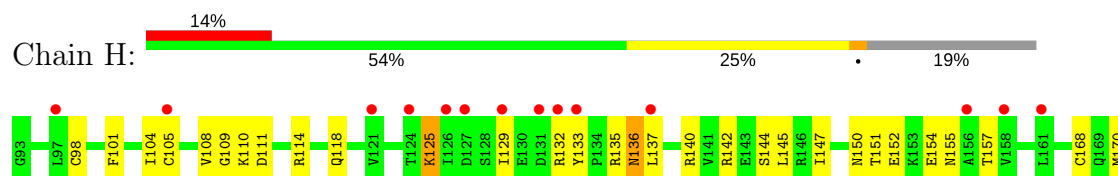


- Molecule 1: Tumor necrosis factor receptor superfamily member 6

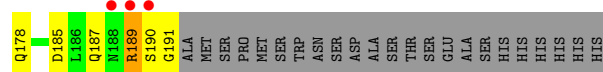
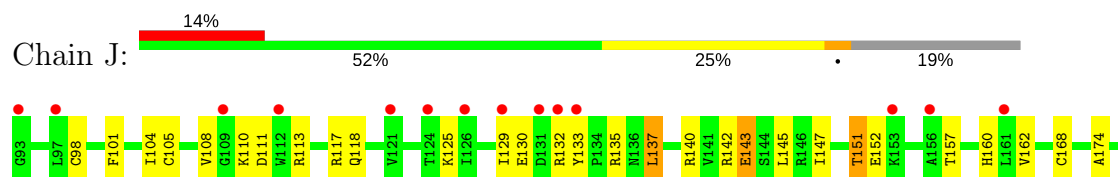




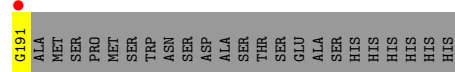
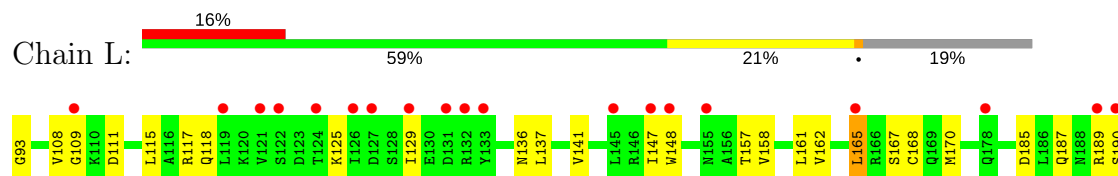
• Molecule 2: Protein FADD



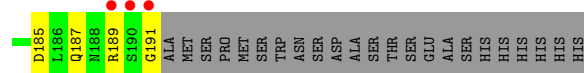
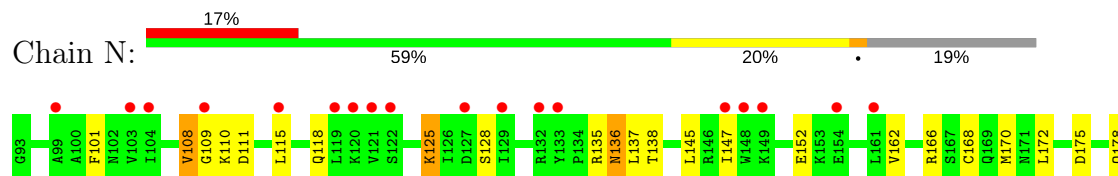
• Molecule 2: Protein FADD



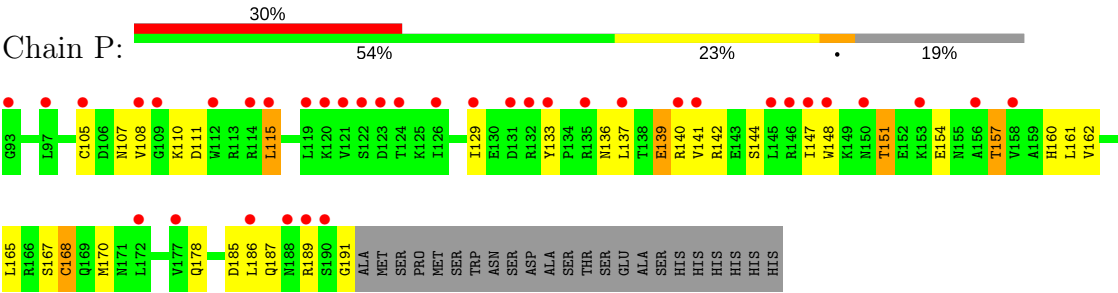
• Molecule 2: Protein FADD



• Molecule 2: Protein FADD



• Molecule 2: Protein FADD



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	126.22Å 126.22Å 299.27Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.73 29.71 – 2.73	Depositor EDS
% Data completeness (in resolution range)	98.8 (30.00-2.73) 98.8 (29.71-2.73)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.234 , 0.278 0.339 , 0.367	Depositor DCC
R_{free} test set	3544 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	73.2	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.499 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13933	wwPDB-VP
Average B, all atoms (Å ²)	87.0	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 45.65 % 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 1.2821e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, 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	0.61	0/923	0.68	0/1241
1	C	0.71	1/923 (0.1%)	0.70	0/1241
1	E	0.67	0/923	0.67	0/1241
1	G	0.73	0/923	0.74	0/1241
1	I	0.70	0/923	0.74	0/1241
1	K	0.64	0/923	0.71	0/1241
1	M	0.72	1/923 (0.1%)	0.72	0/1241
1	O	0.61	0/923	0.66	0/1241
2	B	0.57	0/802	0.61	0/1080
2	D	0.54	1/802 (0.1%)	0.61	0/1080
2	F	0.49	0/802	0.62	0/1080
2	H	0.57	1/802 (0.1%)	0.66	0/1080
2	J	0.55	1/802 (0.1%)	0.65	0/1080
2	L	0.51	0/802	0.59	0/1080
2	N	0.50	0/802	0.63	0/1080
2	P	0.58	2/802 (0.2%)	0.58	0/1080
All	All	0.62	7/13800 (0.1%)	0.66	0/18568

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	279	ARG	NE-CZ	5.70	1.40	1.33
1	M	228	ASP	CG-OD1	5.59	1.38	1.25
2	J	98	CYS	CB-SG	-5.55	1.72	1.81
2	H	98	CYS	CB-SG	-5.33	1.73	1.81
2	D	131	ASP	CG-OD2	5.29	1.37	1.25

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	916	0	939	16	0
1	C	916	0	939	17	0
1	E	916	0	939	14	0
1	G	916	0	939	13	0
1	I	916	0	939	17	0
1	K	916	0	939	21	0
1	M	916	0	939	15	0
1	O	916	0	939	26	0
2	B	795	0	793	24	0
2	D	795	0	793	17	0
2	F	795	0	793	20	0
2	H	795	0	793	20	0
2	J	795	0	793	26	0
2	L	795	0	793	23	0
2	N	795	0	793	19	0
2	P	795	0	793	29	0
3	A	5	0	0	0	0
3	C	5	0	0	0	0
3	E	10	0	0	0	0
3	G	5	0	0	0	0
3	I	5	0	0	0	0
3	J	5	0	0	0	0
3	K	10	0	0	2	0
3	M	10	0	0	1	0
3	O	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	E	1	0	0	0	0
4	H	1	0	0	0	0
4	J	1	0	0	0	0
4	K	1	0	0	0	0
4	M	1	0	0	0	0
4	P	1	0	0	0	0
5	A	7	0	0	1	0
5	B	2	0	0	0	0
5	C	8	0	0	1	0
5	D	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	9	0	0	0	0
5	F	4	0	0	0	0
5	G	11	0	0	0	0
5	H	8	0	0	1	0
5	I	84	0	0	4	0
5	J	8	0	0	0	0
5	K	6	0	0	0	0
5	L	7	0	0	1	0
5	M	9	0	0	1	0
5	N	3	0	0	0	0
5	O	5	0	0	0	0
5	P	2	0	0	3	0
All	All	13933	0	13856	261	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.

The worst 5 of 261 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:L:109:GLY:HA2	2:L:137:LEU:HD21	1.37	1.04
2:B:115:LEU:HD11	2:B:165:LEU:HD23	1.42	1.01
1:K:328:ARG:HE	1:K:332:GLN:HE22	1.13	0.95
2:F:109:GLY:HA2	2:F:137:LEU:HD21	1.47	0.91
1:A:289:GLU:O	1:A:293:THR:HG22	1.73	0.89

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	113/115 (98%)	106 (94%)	6 (5%)	1 (1%)	20	44
1	C	113/115 (98%)	106 (94%)	6 (5%)	1 (1%)	20	44
1	E	113/115 (98%)	106 (94%)	7 (6%)	0	100	100
1	G	113/115 (98%)	108 (96%)	5 (4%)	0	100	100
1	I	113/115 (98%)	108 (96%)	4 (4%)	1 (1%)	20	44
1	K	113/115 (98%)	106 (94%)	6 (5%)	1 (1%)	20	44
1	M	113/115 (98%)	106 (94%)	6 (5%)	1 (1%)	20	44
1	O	113/115 (98%)	106 (94%)	7 (6%)	0	100	100
2	B	97/122 (80%)	92 (95%)	5 (5%)	0	100	100
2	D	97/122 (80%)	92 (95%)	5 (5%)	0	100	100
2	F	97/122 (80%)	93 (96%)	3 (3%)	1 (1%)	18	40
2	H	97/122 (80%)	92 (95%)	5 (5%)	0	100	100
2	J	97/122 (80%)	93 (96%)	4 (4%)	0	100	100
2	L	97/122 (80%)	92 (95%)	5 (5%)	0	100	100
2	N	97/122 (80%)	93 (96%)	3 (3%)	1 (1%)	18	40
2	P	97/122 (80%)	93 (96%)	3 (3%)	1 (1%)	18	40
All	All	1680/1896 (89%)	1592 (95%)	80 (5%)	8 (0%)	32	59

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	256	GLU
1	C	256	GLU
1	K	256	GLU
1	M	256	GLU
1	I	256	GLU

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	105/105 (100%)	103 (98%)	2 (2%)	62	85
1	C	105/105 (100%)	98 (93%)	7 (7%)	19	40
1	E	105/105 (100%)	97 (92%)	8 (8%)	15	33
1	G	105/105 (100%)	99 (94%)	6 (6%)	24	49
1	I	105/105 (100%)	99 (94%)	6 (6%)	24	49
1	K	105/105 (100%)	96 (91%)	9 (9%)	12	26
1	M	105/105 (100%)	99 (94%)	6 (6%)	24	49
1	O	105/105 (100%)	95 (90%)	10 (10%)	10	22
2	B	87/107 (81%)	72 (83%)	15 (17%)	2	5
2	D	87/107 (81%)	83 (95%)	4 (5%)	31	59
2	F	87/107 (81%)	82 (94%)	5 (6%)	24	49
2	H	87/107 (81%)	77 (88%)	10 (12%)	6	14
2	J	87/107 (81%)	77 (88%)	10 (12%)	6	14
2	L	87/107 (81%)	78 (90%)	9 (10%)	8	18
2	N	87/107 (81%)	79 (91%)	8 (9%)	11	23
2	P	87/107 (81%)	74 (85%)	13 (15%)	3	7
All	All	1536/1696 (91%)	1408 (92%)	128 (8%)	13	28

5 of 128 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	239	VAL
2	J	189	ARG
2	P	115	LEU
1	I	270	THR
2	J	117	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 26 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	332	GLN
1	K	280	ASN
2	P	136	ASN
1	G	273	GLN
1	K	252	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

Of 20 ligands modelled in this entry, 8 are monoatomic - leaving 12 for Mogul analysis.

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	A	10	-	4,4,4	0.21	0	6,6,6	0.21	0
3	SO4	C	7	-	4,4,4	0.16	0	6,6,6	0.47	0
3	SO4	E	1	-	4,4,4	0.23	0	6,6,6	0.49	0
3	SO4	E	11	-	4,4,4	0.16	0	6,6,6	0.12	0
3	SO4	G	2	-	4,4,4	0.09	0	6,6,6	0.48	0
3	SO4	I	6	-	4,4,4	0.10	0	6,6,6	0.40	0
3	SO4	J	8	-	4,4,4	0.17	0	6,6,6	0.19	0
3	SO4	K	13	-	4,4,4	0.16	0	6,6,6	0.35	0
3	SO4	K	3	-	4,4,4	0.21	0	6,6,6	0.66	0
3	SO4	M	12	-	4,4,4	0.21	0	6,6,6	0.12	0
3	SO4	M	9	-	4,4,4	0.17	0	6,6,6	0.34	0
3	SO4	O	4	-	4,4,4	0.14	0	6,6,6	0.31	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
3	SO4	A	10	-	-	0/0/0/0	0/0/0/0
3	SO4	C	7	-	-	0/0/0/0	0/0/0/0
3	SO4	E	1	-	-	0/0/0/0	0/0/0/0
3	SO4	E	11	-	-	0/0/0/0	0/0/0/0
3	SO4	G	2	-	-	0/0/0/0	0/0/0/0
3	SO4	I	6	-	-	0/0/0/0	0/0/0/0
3	SO4	J	8	-	-	0/0/0/0	0/0/0/0
3	SO4	K	13	-	-	0/0/0/0	0/0/0/0
3	SO4	K	3	-	-	0/0/0/0	0/0/0/0
3	SO4	M	12	-	-	0/0/0/0	0/0/0/0
3	SO4	M	9	-	-	0/0/0/0	0/0/0/0
3	SO4	O	4	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	K	13	SO4	2	0
3	M	12	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

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	115/115 (100%)	1.05	15 (13%) 4 3	72, 85, 99, 103	0
1	C	115/115 (100%)	1.11	19 (16%) 2 1	72, 85, 99, 103	0
1	E	115/115 (100%)	1.01	15 (13%) 4 3	72, 85, 99, 103	0
1	G	115/115 (100%)	1.25	20 (17%) 2 1	72, 85, 99, 104	0
1	I	115/115 (100%)	1.24	18 (15%) 2 2	72, 85, 99, 104	0
1	K	115/115 (100%)	1.02	15 (13%) 4 3	72, 85, 99, 103	0
1	M	115/115 (100%)	1.10	17 (14%) 3 2	72, 85, 99, 103	0
1	O	115/115 (100%)	1.09	19 (16%) 2 1	72, 85, 99, 103	0
2	B	99/122 (81%)	1.95	31 (31%) 0 0	77, 89, 95, 108	0
2	D	99/122 (81%)	1.27	20 (20%) 1 1	76, 89, 95, 108	0
2	F	99/122 (81%)	1.37	25 (25%) 1 0	76, 89, 95, 108	0
2	H	99/122 (81%)	1.20	17 (17%) 2 1	77, 89, 96, 108	0
2	J	99/122 (81%)	1.18	17 (17%) 2 1	77, 89, 96, 108	0
2	L	99/122 (81%)	1.30	20 (20%) 1 1	76, 89, 95, 108	0
2	N	99/122 (81%)	1.31	21 (21%) 1 1	77, 89, 95, 108	0
2	P	99/122 (81%)	1.95	37 (37%) 0 0	77, 89, 95, 108	0
All	All	1712/1896 (90%)	1.26	326 (19%) 1 1	72, 88, 99, 108	0

The worst 5 of 326 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	P	129	ILE	12.4
2	B	129	ILE	11.1
2	F	129	ILE	10.3
2	J	129	ILE	8.8
2	N	129	ILE	8.8

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	NA	K	16	1/1	0.95	0.49	11.50	62,62,62,62	0
4	NA	M	17	1/1	0.97	0.39	4.63	64,64,64,64	0
4	NA	P	18	1/1	0.33	0.39	4.03	88,88,88,88	0
4	NA	B	21	1/1	0.70	0.36	3.77	84,84,84,84	0
4	NA	E	14	1/1	0.93	0.44	3.61	63,63,63,63	0
4	NA	A	15	1/1	0.96	0.35	1.96	62,62,62,62	0
4	NA	H	20	1/1	0.78	0.32	0.50	77,77,77,77	0
4	NA	J	19	1/1	0.79	0.29	0.17	76,76,76,76	0
3	SO4	M	12	5/5	0.91	0.22	-0.38	117,117,117,118	0
3	SO4	K	3	5/5	0.95	0.19	-0.87	92,95,97,99	0
3	SO4	I	6	5/5	0.92	0.20	-1.38	91,94,96,97	0
3	SO4	E	1	5/5	0.96	0.15	-1.65	95,96,99,101	0
3	SO4	M	9	5/5	0.87	0.17	-1.67	111,111,114,116	0
3	SO4	K	13	5/5	0.93	0.15	-1.98	109,109,110,110	0
3	SO4	A	10	5/5	0.91	0.16	-	114,114,117,117	0
3	SO4	E	11	5/5	0.85	0.21	-	135,135,135,136	0
3	SO4	J	8	5/5	0.93	0.17	-	127,127,128,128	0
3	SO4	C	7	5/5	0.84	0.24	-	120,120,121,124	0
3	SO4	G	2	5/5	0.89	0.21	-	95,95,97,98	0
3	SO4	O	4	5/5	0.88	0.19	-	116,117,119,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.