



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:19 am GMT

PDB ID : 4KXZ  
Title : crystal structure of tgfb2 in complex with GC2008.  
Authors : Mathieu, M.; Moulin, A.G.; Wei, R.  
Deposited on : 2013-05-28  
Resolution : 2.83 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](mailto: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.

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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

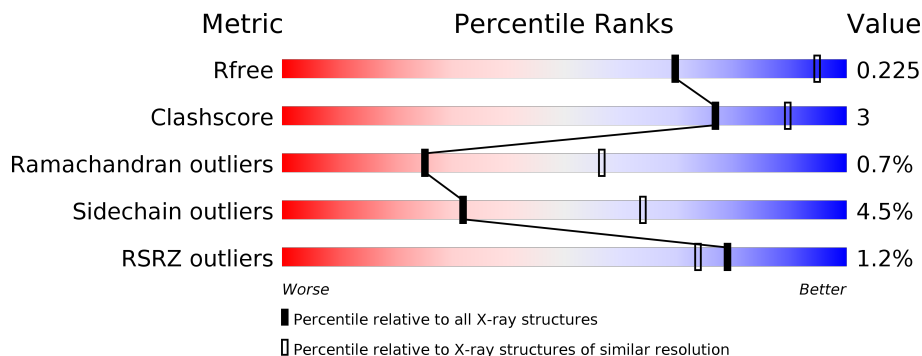
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.83 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	3466 (2.88-2.80)
Clashscore	112137	3975 (2.88-2.80)
Ramachandran outliers	110173	3902 (2.88-2.80)
Sidechain outliers	110143	3905 (2.88-2.80)
RSRZ outliers	101464	3501 (2.88-2.80)

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  $\geq 3$ , 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	112	<div> <div>88%</div> <div>12%</div> </div>
1	B	112	<div>4%</div> <div>89%</div> <div>11%</div>
1	D	112	<div>4%</div> <div>89%</div> <div>11%</div>
1	E	112	<div>88%</div> <div>13%</div>
2	H	224	<div>%</div> <div>86%</div> <div>13%</div> <div>.</div>
2	J	224	<div>2%</div> <div>80%</div> <div>16%</div> <div>..</div>

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Mol	Chain	Length	Quality of chain
2	N	224	
2	Q	224	
3	I	215	
3	L	215	
3	M	215	
3	P	215	

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	MES	E	201	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16951 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 Transforming growth factor beta-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	112	Total	C	N	O	S	0	0	0
			890	566	149	165	10			
1	B	112	Total	C	N	O	S	0	0	0
			890	566	149	165	10			
1	D	112	Total	C	N	O	S	0	0	0
			890	566	149	165	10			
1	E	112	Total	C	N	O	S	0	0	0
			890	566	149	165	10			

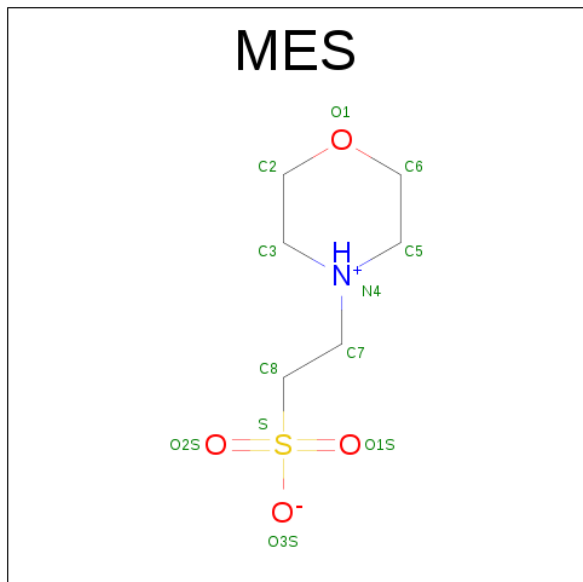
- Molecule 2 is a protein called GC1008 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	223	Total	C	N	O	S	0	0	0
			1670	1048	284	330	8			
2	J	217	Total	C	N	O	S	0	0	0
			1615	1017	269	321	8			
2	N	223	Total	C	N	O	S	0	0	0
			1670	1048	284	330	8			
2	Q	217	Total	C	N	O	S	0	0	0
			1618	1019	270	321	8			

- Molecule 3 is a protein called GC1008 Light Chain.

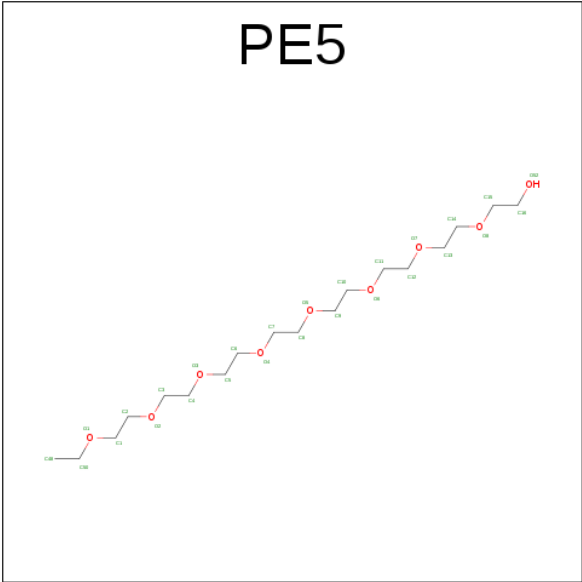
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	212	Total	C	N	O	S	0	0	0
			1623	1015	277	327	4			
3	L	213	Total	C	N	O	S	0	0	0
			1632	1020	278	330	4			
3	M	213	Total	C	N	O	S	0	0	0
			1632	1020	278	330	4			
3	P	213	Total	C	N	O	S	0	0	0
			1632	1020	278	330	4			

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula:  $C_6H_{13}NO_4S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	E	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	J	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	Q	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	M	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is 3,6,9,12,15,18,21,24-OCTAOXAHEXACOSAN-1-OL (three-letter code: PE5) (formula:  $C_{18}H_{38}O_9$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	P	1	Total	C	O	0	0
			13	9	4		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	17	Total	O	0	0
			17	17		
6	B	5	Total	O	0	0
			5	5		
6	D	12	Total	O	0	0
			12	12		
6	E	14	Total	O	0	0
			14	14		
6	H	31	Total	O	0	0
			31	31		
6	J	17	Total	O	0	0
			17	17		
6	N	28	Total	O	0	0
			28	28		
6	Q	16	Total	O	0	0
			16	16		
6	I	7	Total	O	0	0
			7	7		
6	L	30	Total	O	0	0
			30	30		
6	M	35	Total	O	0	0
			35	35		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	P	14	Total	O	0	0
			14	14		

### 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: Transforming growth factor beta-2

Chain A: 




- Molecule 1: Transforming growth factor beta-2

Chain B: 




- Molecule 1: Transforming growth factor beta-2

Chain D: 




- Molecule 1: Transforming growth factor beta-2

Chain E: 



- Molecule 2: GC1008 Heavy Chain

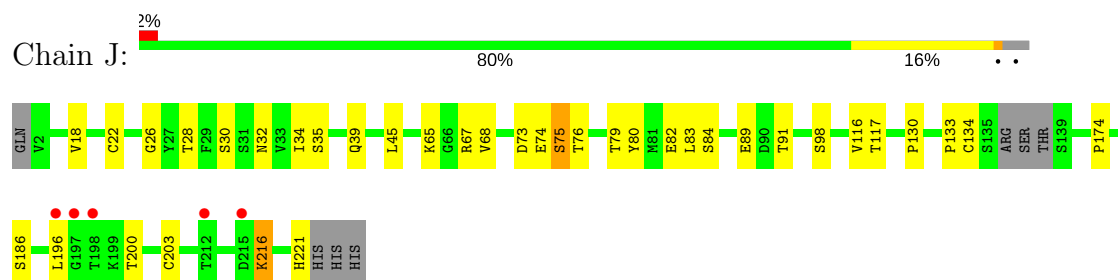
Chain H: 



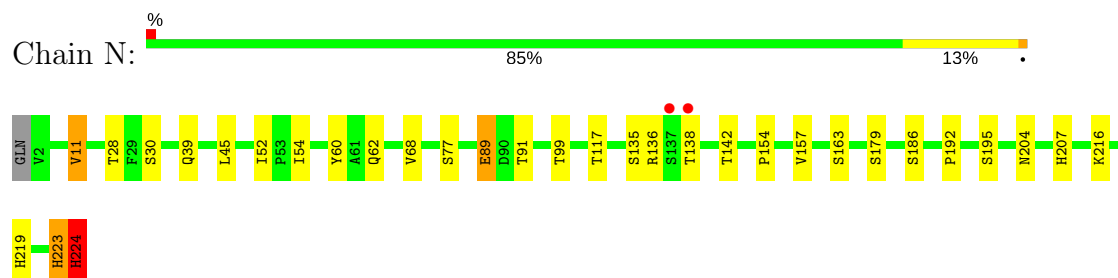
H224



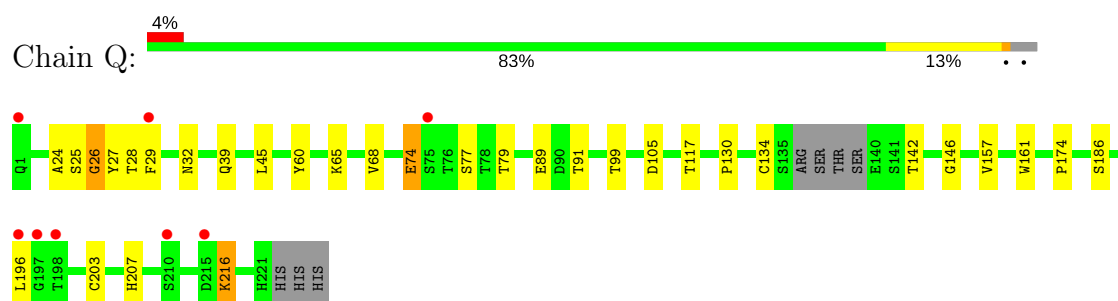
## • Molecule 2: GC1008 Heavy Chain



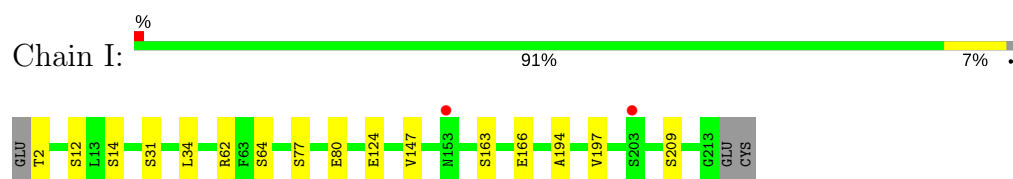
## • Molecule 2: GC1008 Heavy Chain



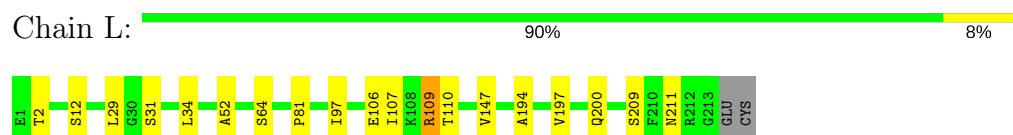
## • Molecule 2: GC1008 Heavy Chain



## • Molecule 3: GC1008 Light Chain



## • Molecule 3: GC1008 Light Chain

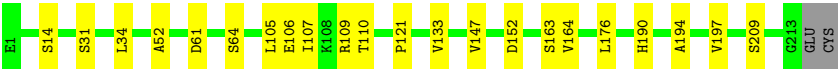


## • Molecule 3: GC1008 Light Chain





● Molecule 3: GC1008 Light Chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	131.20Å 359.68Å 64.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.73 – 2.83 48.73 – 2.83	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.73-2.83) 99.5 (48.73-2.83)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 2.81Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, $R_{free}$	0.183 , 0.222 0.189 , 0.225	Depositor DCC
$R_{free}$ test set	3709 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	52.3	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 49.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	16951	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PE5, MES

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	0/914	0.70	0/1242
1	B	0.50	0/914	0.71	0/1242
1	D	0.52	0/914	0.69	0/1242
1	E	0.50	0/914	0.68	0/1242
2	H	0.56	0/1710	0.78	1/2332 (0.0%)
2	J	0.49	0/1651	0.76	0/2252
2	N	0.54	0/1710	0.77	1/2332 (0.0%)
2	Q	0.51	0/1654	0.75	0/2256
3	I	0.46	0/1659	0.68	0/2254
3	L	0.51	0/1668	0.71	0/2266
3	M	0.53	0/1668	0.71	0/2266
3	P	0.48	0/1668	0.69	0/2266
All	All	0.51	0/17044	0.72	2/23192 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	224	HIS	CB-CA-C	6.17	122.73	110.40
2	N	224	HIS	CB-CA-C	5.42	121.23	110.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	890	0	860	10	0
1	B	890	0	860	10	0
1	D	890	0	859	5	0
1	E	890	0	859	6	0
2	H	1670	0	1635	24	0
2	J	1615	0	1588	15	0
2	N	1670	0	1635	21	0
2	Q	1618	0	1594	16	0
3	I	1623	0	1577	6	0
3	L	1632	0	1586	8	0
3	M	1632	0	1586	4	0
3	P	1632	0	1586	7	0
4	A	12	0	13	0	0
4	E	12	0	13	1	0
4	J	12	0	13	0	0
4	M	12	0	13	0	0
4	Q	12	0	13	0	0
5	P	13	0	14	0	0
6	A	17	0	0	0	0
6	B	5	0	0	0	0
6	D	12	0	0	0	0
6	E	14	0	0	0	0
6	H	31	0	0	4	0
6	I	7	0	0	2	0
6	J	17	0	0	0	0
6	L	30	0	0	0	0
6	M	35	0	0	0	0
6	N	28	0	0	4	0
6	P	14	0	0	0	0
6	Q	16	0	0	0	0
All	All	16951	0	16304	114	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 3.

All (114) 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:N:11:VAL:HG22	2:N:154:PRO:HG3	1.51	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:11:VAL:CG2	2:N:154:PRO:HG3	2.18	0.73
1:A:77:CYS:SG	1:B:77:CYS:HB3	2.27	0.73
3:L:81:PRO:HA	3:L:107:ILE:HG21	1.72	0.71
2:H:224:HIS:ND1	6:H:329:HOH:O	2.24	0.68
2:H:219:HIS:ND1	2:H:224:HIS:CG	2.63	0.67
2:N:219:HIS:ND1	2:N:224:HIS:CG	2.63	0.67
2:Q:24:ALA:HB2	2:Q:29:PHE:CE2	2.29	0.66
1:A:36:PRO:HD2	1:A:88:ILE:HG22	1.78	0.66
3:I:80:GLU:OE1	6:I:307:HOH:O	2.14	0.65
2:N:224:HIS:ND1	6:N:327:HOH:O	2.29	0.64
2:H:11:VAL:HG22	2:H:154:PRO:HG3	1.80	0.63
2:H:104:LEU:HG	3:L:97:ILE:HG13	1.80	0.63
2:N:219:HIS:ND1	2:N:224:HIS:ND1	2.48	0.61
2:N:219:HIS:ND1	2:N:224:HIS:CE1	2.69	0.60
2:N:223:HIS:CD2	6:N:326:HOH:O	2.55	0.58
3:P:164:VAL:HG22	3:P:176:LEU:HD12	1.85	0.58
1:D:83:LEU:HD13	1:D:103:ASN:HB3	1.86	0.58
2:H:219:HIS:ND1	2:H:224:HIS:CE1	2.73	0.57
3:L:109:ARG:HD3	3:L:110:THR:O	2.06	0.56
1:A:26:ARG:HD2	1:A:27:ASP:OD1	2.05	0.56
2:J:34:ILE:HG21	2:J:79:THR:HG21	1.89	0.55
1:E:83:LEU:HD13	1:E:103:ASN:HB3	1.87	0.55
2:H:224:HIS:HD2	2:H:224:HIS:OXT	1.90	0.55
1:A:83:LEU:HD13	1:A:103:ASN:HB3	1.87	0.55
2:J:67:ARG:HD2	2:J:84:SER:O	2.06	0.55
2:H:219:HIS:ND1	2:H:224:HIS:ND1	2.55	0.54
3:I:2:THR:N	6:I:302:HOH:O	2.40	0.54
2:Q:39:GLN:HB2	2:Q:45:LEU:HD23	1.90	0.54
1:A:50:TYR:CZ	2:J:30:SER:HB3	2.43	0.53
2:N:39:GLN:HB2	2:N:45:LEU:HD23	1.90	0.52
2:H:13:LYS:NZ	6:H:305:HOH:O	2.24	0.51
2:J:39:GLN:HB2	2:J:45:LEU:HD23	1.92	0.51
2:N:224:HIS:OXT	2:N:224:HIS:HD2	1.93	0.51
3:P:109:ARG:NH1	3:P:110:THR:O	2.43	0.51
1:B:60:ARG:HH21	2:H:54:ILE:HA	1.76	0.50
1:B:26:ARG:HD2	1:B:27:ASP:OD1	2.11	0.50
3:M:194:ALA:HB2	3:M:209:SER:HB3	1.94	0.50
2:H:11:VAL:CG2	2:H:154:PRO:HG3	2.41	0.50
3:L:194:ALA:HB2	3:L:209:SER:HB3	1.93	0.50
2:N:157:VAL:HG12	2:N:207:HIS:HB2	1.93	0.49
1:A:77:CYS:SG	1:B:77:CYS:CB	3.00	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:LEU:HD23	1:D:106:VAL:HA	1.94	0.49
1:E:26:ARG:HD2	1:E:27:ASP:OD1	2.12	0.49
2:H:219:HIS:ND1	2:H:224:HIS:CD2	2.80	0.49
2:J:28:THR:O	2:J:32:ASN:HB2	2.13	0.49
2:Q:74:GLU:CD	2:Q:74:GLU:H	2.16	0.48
2:H:64:PHE:HB3	2:H:68:VAL:CG2	2.43	0.48
2:N:223:HIS:NE2	6:N:326:HOH:O	2.34	0.48
2:H:224:HIS:CE1	6:H:329:HOH:O	2.62	0.48
2:N:224:HIS:OXT	2:N:224:HIS:CD2	2.67	0.48
3:L:147:VAL:HG22	3:L:197:VAL:HG22	1.95	0.48
2:H:224:HIS:OXT	2:H:224:HIS:CD2	2.66	0.48
1:D:79:VAL:HG11	1:E:79:VAL:HG11	1.95	0.48
3:P:147:VAL:HG22	3:P:197:VAL:HG22	1.95	0.48
3:I:62:ARG:HD2	3:I:77:SER:O	2.14	0.47
2:Q:130:PRO:HD3	2:Q:216:LYS:HE2	1.96	0.47
1:A:77:CYS:SG	1:B:77:CYS:SG	3.11	0.47
2:Q:28:THR:O	2:Q:32:ASN:HB2	2.14	0.47
2:H:223:HIS:ND1	6:H:328:HOH:O	2.14	0.46
1:D:39:TYR:CE2	1:D:41:ALA:HB2	2.51	0.46
2:J:73:ASP:HB2	2:J:80:TYR:HE2	1.81	0.46
2:Q:174:PRO:HD2	3:P:163:SER:OG	2.16	0.46
2:H:68:VAL:HG22	2:H:83:LEU:HD13	1.98	0.46
3:I:194:ALA:HB2	3:I:209:SER:HB3	1.97	0.46
4:E:201:MES:H72	2:H:14:PRO:O	2.15	0.46
2:N:224:HIS:CE1	6:N:327:HOH:O	2.68	0.46
1:D:50:TYR:OH	2:N:30:SER:HB3	2.15	0.46
2:H:52:ILE:HD12	2:H:57:ILE:HB	1.97	0.46
3:I:147:VAL:HG22	3:I:197:VAL:HG22	1.98	0.46
2:J:74:GLU:HG2	2:J:75:SER:H	1.80	0.46
3:M:147:VAL:HG22	3:M:197:VAL:HG22	1.98	0.46
2:J:68:VAL:HG22	2:J:83:LEU:HD13	1.98	0.46
2:Q:157:VAL:HG12	2:Q:207:HIS:HB2	1.98	0.46
3:P:194:ALA:HB2	3:P:209:SER:HB3	1.98	0.45
1:B:102:SER:HB3	2:N:163:SER:OG	2.17	0.45
2:J:91:THR:HG23	2:J:117:THR:HA	1.98	0.45
1:E:32:TRP:CE2	2:N:52:ILE:HG21	2.52	0.44
2:N:91:THR:HG23	2:N:117:THR:HA	1.99	0.44
2:J:174:PRO:HD2	3:I:163:SER:OG	2.17	0.44
3:P:152:ASP:OD1	3:P:190:HIS:HB3	2.17	0.44
2:N:60:TYR:CD2	2:N:68:VAL:HG13	2.53	0.44
2:Q:60:TYR:HB2	2:Q:65:LYS:HG3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:219:HIS:ND1	2:N:224:HIS:CD2	2.87	0.43
2:Q:29:PHE:CE1	2:Q:79:THR:HG23	2.54	0.43
2:Q:91:THR:HG23	2:Q:117:THR:HA	2.01	0.43
1:E:83:LEU:HD23	1:E:106:VAL:HA	2.01	0.43
2:J:91:THR:HA	2:J:116:VAL:O	2.18	0.43
2:Q:26:GLY:C	2:Q:28:THR:H	2.22	0.43
1:E:67:THR:HG21	2:Q:27:TYR:HE1	1.83	0.43
3:L:2:THR:HG21	3:L:29:LEU:HD21	2.01	0.42
2:Q:146:GLY:HA2	2:Q:161:TRP:CH2	2.55	0.42
2:J:130:PRO:HD3	2:J:216:LYS:HE2	2.01	0.42
3:M:62:ARG:NH1	3:M:83:ASP:OD2	2.50	0.42
1:B:50:TYR:OH	2:H:30:SER:HB3	2.19	0.42
2:H:64:PHE:HB3	2:H:68:VAL:HG21	2.00	0.41
3:L:12:SER:HA	3:L:106:GLU:O	2.19	0.41
2:H:198:THR:HG22	2:Q:65:LYS:NZ	2.35	0.41
2:J:74:GLU:HG2	2:J:75:SER:N	2.36	0.41
2:N:89:GLU:H	2:N:89:GLU:CD	2.24	0.41
2:Q:99:THR:HB	2:Q:105:ASP:HB3	2.02	0.41
1:B:50:TYR:HA	1:B:63:SER:HB2	2.02	0.41
2:Q:60:TYR:CD2	2:Q:68:VAL:HG13	2.56	0.41
2:J:18:VAL:O	2:J:82:GLU:HA	2.21	0.41
2:N:192:PRO:O	2:N:195:SER:HB2	2.20	0.41
1:A:32:TRP:CE2	2:H:52:ILE:HG21	2.56	0.40
2:H:136:ARG:NH1	3:L:211:ASN:H	2.19	0.40
3:P:121:PRO:HD3	3:P:133:VAL:HG22	2.02	0.40
1:A:74:ALA:HA	1:B:43:PHE:CD2	2.56	0.40
1:B:23:ASP:HB3	1:B:26:ARG:HB3	2.02	0.40
1:A:34:HIS:CD2	1:A:91:TYR:HE1	2.39	0.40
2:J:22:CYS:HB3	2:J:79:THR:HG23	2.02	0.40
3:M:62:ARG:HH11	3:M:83:ASP:CG	2.24	0.40
2:H:192:PRO:O	2:H:195:SER:HB2	2.21	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	110/112 (98%)	105 (96%)	5 (4%)	0	100	100
1	B	110/112 (98%)	103 (94%)	7 (6%)	0	100	100
1	D	110/112 (98%)	105 (96%)	5 (4%)	0	100	100
1	E	110/112 (98%)	103 (94%)	7 (6%)	0	100	100
2	H	221/224 (99%)	210 (95%)	8 (4%)	3 (1%)	13	37
2	J	213/224 (95%)	201 (94%)	8 (4%)	4 (2%)	9	28
2	N	221/224 (99%)	209 (95%)	10 (4%)	2 (1%)	20	49
2	Q	213/224 (95%)	202 (95%)	8 (4%)	3 (1%)	13	37
3	I	210/215 (98%)	201 (96%)	9 (4%)	0	100	100
3	L	211/215 (98%)	198 (94%)	11 (5%)	2 (1%)	20	49
3	M	211/215 (98%)	200 (95%)	10 (5%)	1 (0%)	32	64
3	P	211/215 (98%)	201 (95%)	9 (4%)	1 (0%)	32	64
All	All	2151/2204 (98%)	2038 (95%)	97 (4%)	16 (1%)	25	56

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	N	135	SER
2	H	223	HIS
2	J	26	GLY
2	J	75	SER
2	Q	26	GLY
2	Q	196	LEU
3	L	200	GLN
2	H	42	GLY
2	H	139	SER
2	J	76	THR
2	N	223	HIS
3	L	52	ALA
3	M	52	ALA
2	Q	25	SER
3	P	52	ALA
2	J	133	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	100/100 (100%)	99 (99%)	1 (1%)	80	94
1	B	100/100 (100%)	97 (97%)	3 (3%)	46	78
1	D	100/100 (100%)	95 (95%)	5 (5%)	28	59
1	E	100/100 (100%)	94 (94%)	6 (6%)	22	51
2	H	191/192 (100%)	179 (94%)	12 (6%)	21	49
2	J	185/192 (96%)	174 (94%)	11 (6%)	23	52
2	N	191/192 (100%)	176 (92%)	15 (8%)	14	37
2	Q	185/192 (96%)	177 (96%)	8 (4%)	33	66
3	I	182/185 (98%)	175 (96%)	7 (4%)	38	70
3	L	183/185 (99%)	179 (98%)	4 (2%)	57	85
3	M	183/185 (99%)	178 (97%)	5 (3%)	50	81
3	P	183/185 (99%)	175 (96%)	8 (4%)	33	65
All	All	1883/1908 (99%)	1798 (96%)	85 (4%)	32	64

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	CYS
1	B	16	CYS
1	B	53	SER
1	B	71	GLU
1	D	16	CYS
1	D	31	LYS
1	D	77	CYS
1	D	80	SER
1	D	110	LYS
1	E	16	CYS
1	E	31	LYS
1	E	53	SER
1	E	59	SER
1	E	77	CYS

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Mol	Chain	Res	Type
1	E	89	LEU
2	H	11	VAL
2	H	31	SER
2	H	35	SER
2	H	62	GLN
2	H	77	SER
2	H	89	GLU
2	H	138	THR
2	H	142	THR
2	H	186	SER
2	H	216	LYS
2	H	217	ARG
2	H	224	HIS
2	J	35	SER
2	J	65	LYS
2	J	89	GLU
2	J	98	SER
2	J	134	CYS
2	J	186	SER
2	J	196	LEU
2	J	200	THR
2	J	203	CYS
2	J	216	LYS
2	J	221	HIS
2	N	11	VAL
2	N	28	THR
2	N	54	ILE
2	N	62	GLN
2	N	77	SER
2	N	89	GLU
2	N	99	THR
2	N	136	ARG
2	N	138	THR
2	N	142	THR
2	N	179	SER
2	N	186	SER
2	N	204	ASN
2	N	216	LYS
2	N	224	HIS
2	Q	74	GLU
2	Q	77	SER
2	Q	89	GLU

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Mol	Chain	Res	Type
2	Q	134	CYS
2	Q	142	THR
2	Q	186	SER
2	Q	203	CYS
2	Q	216	LYS
3	I	12	SER
3	I	14	SER
3	I	31	SER
3	I	34	LEU
3	I	64	SER
3	I	124	GLU
3	I	166	GLU
3	L	31	SER
3	L	34	LEU
3	L	64	SER
3	L	109	ARG
3	M	14	SER
3	M	31	SER
3	M	34	LEU
3	M	62	ARG
3	M	64	SER
3	P	14	SER
3	P	31	SER
3	P	34	LEU
3	P	61	ASP
3	P	64	SER
3	P	105	LEU
3	P	106	GLU
3	P	107	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

6 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	MES	A	201	-	12,12,12	0.87	1 (8%)	14,16,16	0.31	0
4	MES	E	201	-	12,12,12	0.70	0	14,16,16	0.33	0
4	MES	J	301	-	12,12,12	1.49	1 (8%)	14,16,16	0.34	0
4	MES	M	301	-	12,12,12	0.59	0	14,16,16	0.45	0
5	PE5	P	701	-	12,12,26	0.21	0	11,11,25	0.20	0
4	MES	Q	801	-	12,12,12	1.26	1 (8%)	14,16,16	0.38	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MES	A	201	-	-	0/6/14/14	0/1/1/1
4	MES	E	201	-	-	0/6/14/14	0/1/1/1
4	MES	J	301	-	-	0/6/14/14	0/1/1/1
4	MES	M	301	-	-	0/6/14/14	0/1/1/1
5	PE5	P	701	-	-	0/10/10/24	0/0/0/0
4	MES	Q	801	-	-	0/6/14/14	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	J	301	MES	C8-S	-4.78	1.70	1.77
4	Q	801	MES	C8-S	-3.92	1.71	1.77
4	A	201	MES	C8-S	-2.13	1.74	1.77

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	201	MES	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	112/112 (100%)	-0.48	0 100 100	27, 43, 59, 73	0
1	B	112/112 (100%)	-0.05	4 (3%) 43 33	32, 47, 70, 88	0
1	D	112/112 (100%)	-0.01	4 (3%) 43 33	34, 52, 77, 103	0
1	E	112/112 (100%)	-0.25	0 100 100	31, 47, 65, 91	0
2	H	223/224 (99%)	-0.33	2 (0%) 84 80	27, 41, 73, 101	0
2	J	217/224 (96%)	0.06	5 (2%) 61 52	33, 57, 89, 107	0
2	N	223/224 (99%)	-0.36	2 (0%) 84 80	28, 46, 84, 99	0
2	Q	217/224 (96%)	0.02	8 (3%) 42 32	26, 52, 95, 114	0
3	I	212/215 (98%)	-0.09	2 (0%) 84 80	35, 56, 79, 86	0
3	L	213/215 (99%)	-0.10	0 100 100	26, 47, 72, 86	0
3	M	213/215 (99%)	-0.33	0 100 100	25, 42, 69, 82	0
3	P	213/215 (99%)	-0.28	0 100 100	26, 47, 77, 89	0
All	All	2179/2204 (98%)	-0.18	27 (1%) 79 73	25, 48, 79, 114	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	Q	1	GLN	4.7
2	J	196	LEU	4.5
2	J	197	GLY	3.4
1	D	112	SER	3.4
2	J	212	THR	3.0
2	N	138	THR	2.9
2	Q	29	PHE	2.9
2	Q	196	LEU	2.7
1	D	7	CYS	2.5
3	I	153	ASN	2.5
1	B	6	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
3	I	203	SER	2.4
2	J	215	ASP	2.4
2	Q	75	SER	2.3
2	Q	198	THR	2.3
2	J	198	THR	2.3
2	N	137	SER	2.2
2	Q	215	ASP	2.2
1	D	9	ARG	2.1
2	Q	210	SER	2.1
1	B	10	ASN	2.1
2	H	27	TYR	2.1
1	B	11	VAL	2.1
1	B	9	ARG	2.1
1	D	71	GLU	2.0
2	Q	197	GLY	2.0
2	H	136	ARG	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	MES	E	201	12/12	0.80	0.34	5.55	129,131,132,132	0
4	MES	M	301	12/12	0.87	0.21	1.67	63,75,103,103	0
5	PE5	P	701	13/27	0.94	0.23	1.39	48,52,65,67	0
4	MES	J	301	12/12	0.95	0.26	0.93	70,72,74,77	0
4	MES	Q	801	12/12	0.93	0.25	0.58	62,65,72,75	0
4	MES	A	201	12/12	0.96	0.18	0.45	57,61,66,67	0



## 6.5 Other polymers [i](#)

There are no such residues in this entry.