



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:47 PM GMT

PDB ID : 3V0A  
Title : 2.7 angstrom crystal structure of BoNT/Ai in complex with NTNHA  
Authors : Gu, S.; Rumpel, S.; Zhou, J.; Strotmeier, J.; Bigalke, H.; Perry, K.; Shoemaker, C.B.; Rummel, A.; Jin, R.  
Deposited on : 2011-12-07  
Resolution : 2.70 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

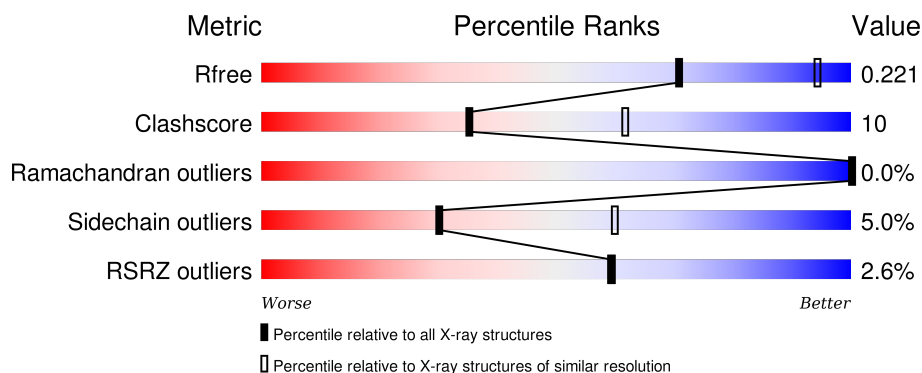
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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  $\geq 3$ , 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	1296	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 80%, yellow 80%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>80%</span> <span>17%</span> </div> </div>
2	B	1196	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 2%, green 2%, green 73%, yellow 73%, yellow 94%, grey 94%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>73%</span> <span>21%</span> </div> </div>
3	C	152	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 18%, green 18%, green 43%, yellow 43%, yellow 75%, grey 75%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> <span>43%</span> <span>32%</span> <span>22%</span> </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
6	SO4	A	1301	-	-	-	X
6	SO4	A	1303	-	-	-	X
6	SO4	A	1304	-	-	-	X
6	SO4	A	1307	-	-	-	X
6	SO4	A	1309	-	-	-	X
6	SO4	A	1310	-	-	-	X
6	SO4	B	1196	-	-	-	X
6	SO4	B	1199	-	-	-	X
6	SO4	B	1201	-	-	-	X
7	MES	B	1204	-	-	-	X
7	MES	B	1205	-	-	-	X

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 21163 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 BoNT/A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1280	Total	C	N	O	S	36	0	0
			10418	6681	1721	1983	33			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	224	GLN	GLU	ENGINEERED MUTATION	UNP Q7B8V4
A	363	ALA	ARG	ENGINEERED MUTATION	UNP Q7B8V4
A	366	PHE	TYR	ENGINEERED MUTATION	UNP Q7B8V4
A	1158	ALA	THR	CONFLICT	UNP Q7B8V4

- Molecule 2 is a protein called NTN1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	1150	Total	C	N	O	S	46	0	0
			9409	6022	1515	1840	32			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	EXPRESSION TAG	UNP Q45914
B	0	SER	-	EXPRESSION TAG	UNP Q45914
B	1194	PRO	-	EXPRESSION TAG	UNP Q45914

- Molecule 3 is a protein called Llama antibody F12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	119	Total	C	N	O	S	4	0	0
			892	551	156	181	4			

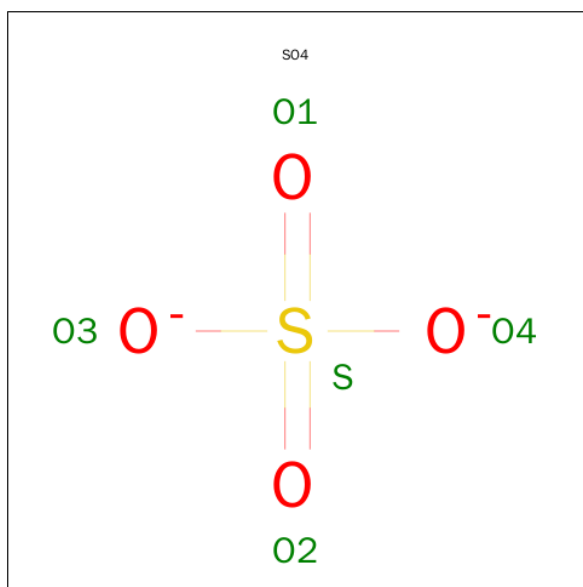
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



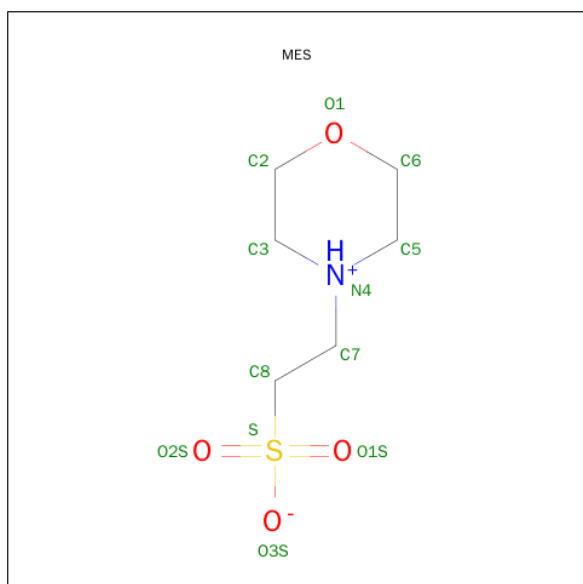
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 7 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>S).



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

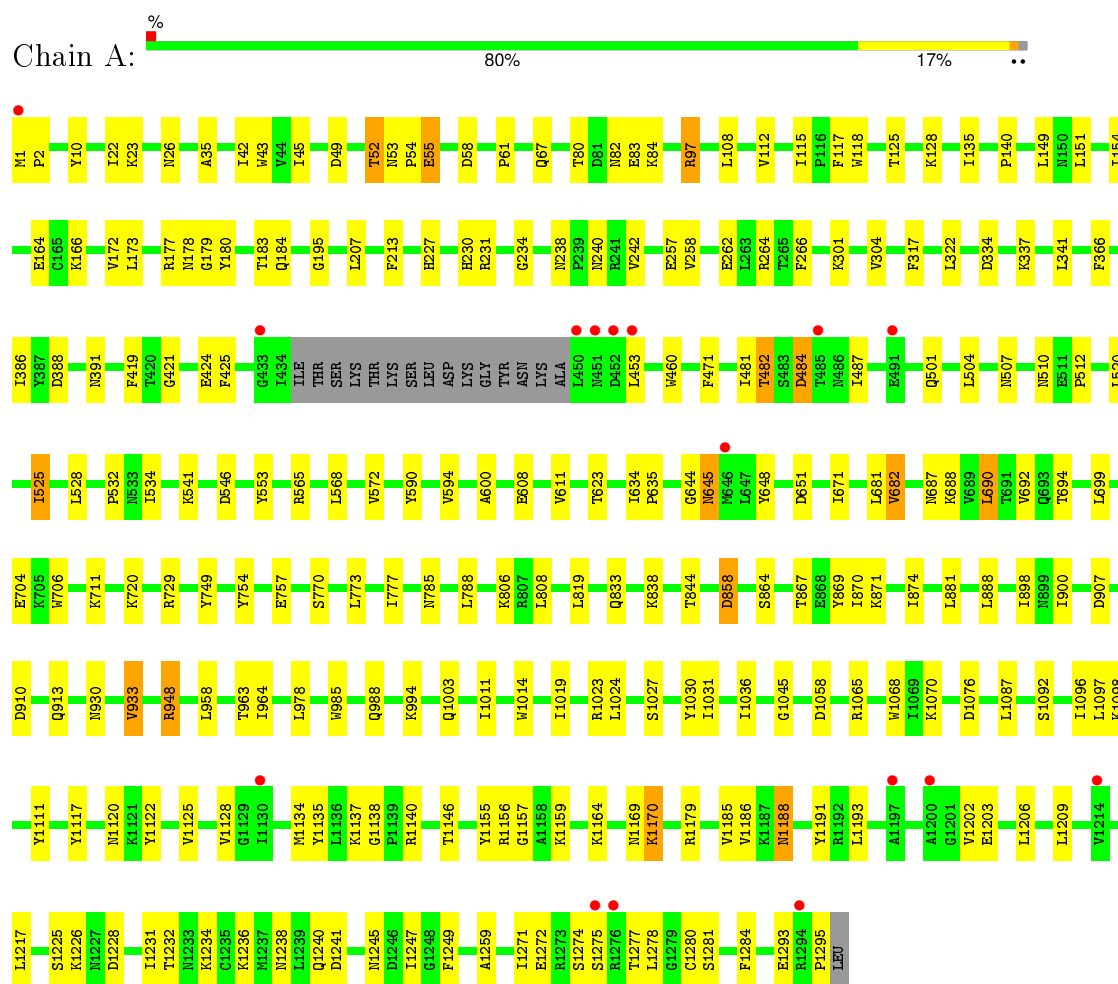
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	181	Total	O	0	0
			181	181		
8	B	112	Total	O	0	0
			112	112		
8	C	1	Total	O	0	0
			1	1		

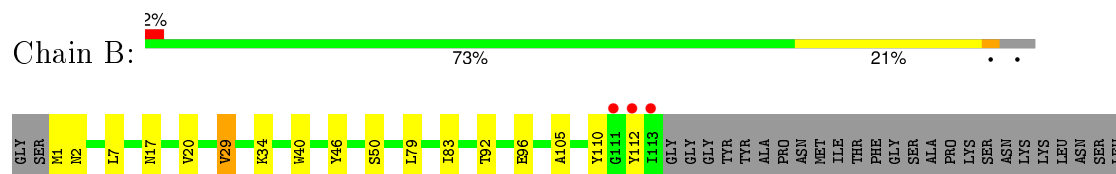
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

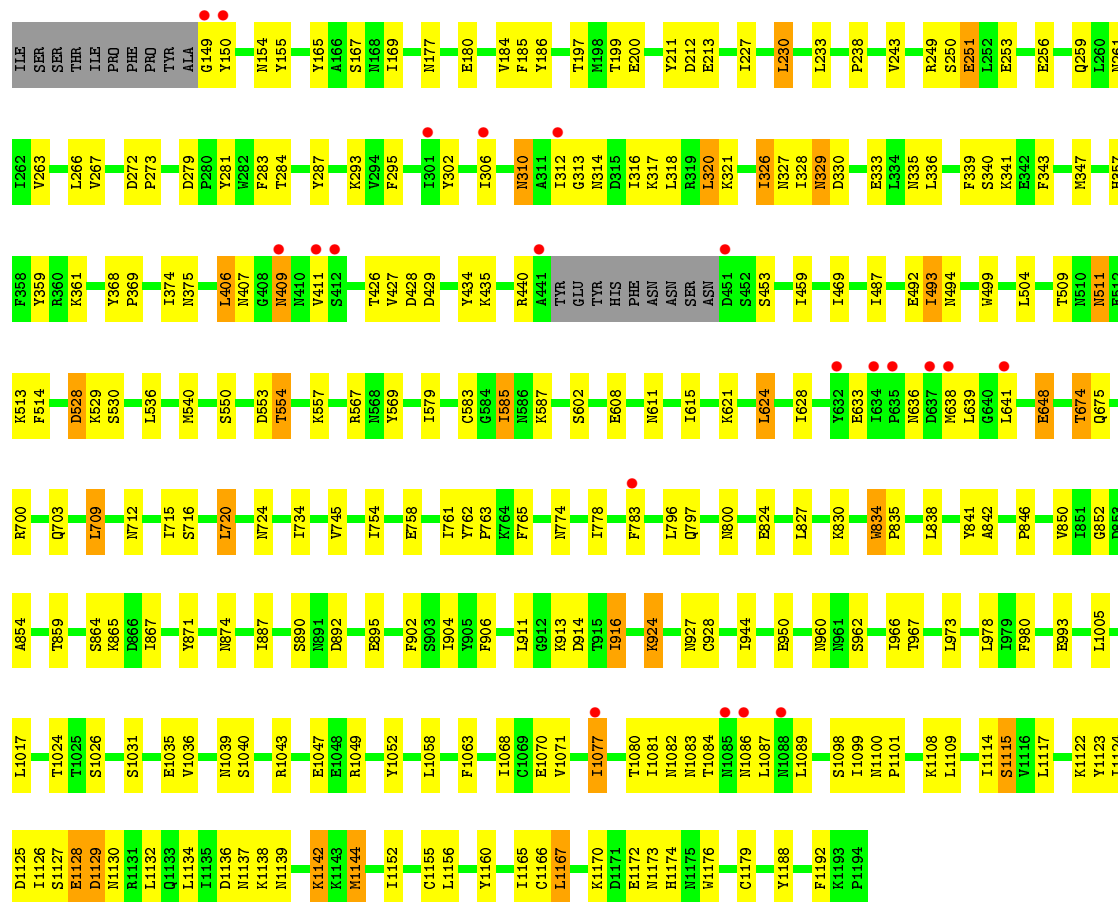
#### • Molecule 1: BoNT/A



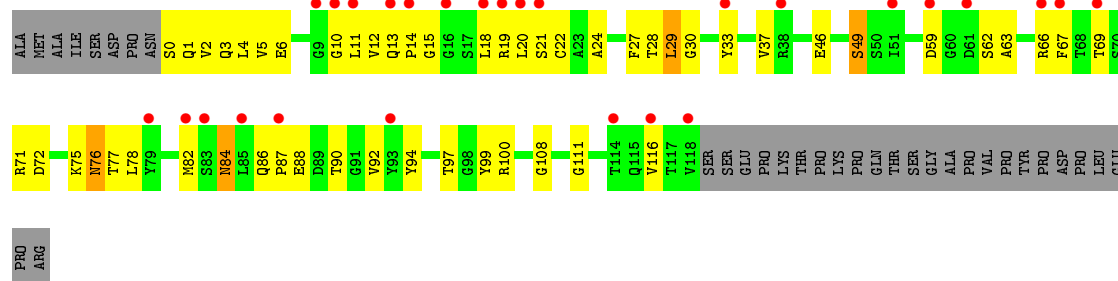
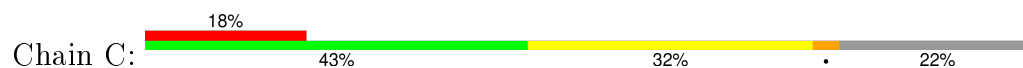
#### • Molecule 2: NTNH







### • Molecule 3: Llama antibody F12



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	156.08 Å   156.08 Å   324.60 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	49.36 – 2.70 49.54 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.36-2.70) 99.7 (49.54-2.70)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.89 (at 2.69 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, $R_{free}$	0.176   ,   0.224 0.171   ,   0.221	Depositor DCC
$R_{free}$ test set	5467 reflections (4.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.5	Xtriage
Anisotropy	0.556	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 49.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 109641 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	21163	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.96% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN, CA, MES, 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	0.31	0/10639	0.49	0/14405
2	B	0.32	0/9604	0.51	0/13015
3	C	0.24	0/912	0.45	0/1238
All	All	0.31	0/21155	0.50	0/28658

There are no bond length outliers.

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	10418	0	10297	180	0
2	B	9409	0	9202	194	0
3	C	892	0	840	41	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
6	A	60	0	0	2	0
6	B	35	0	0	1	0
6	C	5	0	0	0	0
7	B	48	0	52	1	0
8	A	181	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	112	0	0	2	0
8	C	1	0	0	1	0
All	All	21163	0	20391	404	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 10.

All (404) 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:A:1275:SER:HA	2:B:827:LEU:HD21	1.22	1.21
2:B:528:ASP:HB2	2:B:530:SER:H	1.07	1.13
2:B:528:ASP:HB2	2:B:530:SER:N	1.76	1.00
2:B:426:THR:HB	2:B:427:VAL:CA	1.91	0.99
2:B:426:THR:HB	2:B:427:VAL:HA	1.02	0.98
2:B:426:THR:CB	2:B:427:VAL:HA	1.93	0.98
1:A:421:GLY:H	1:A:424:GLU:HG3	1.33	0.93
1:A:1275:SER:CA	2:B:827:LEU:HD21	2.04	0.87
2:B:528:ASP:CB	2:B:530:SER:H	1.88	0.84
2:B:314:ASN:HB3	2:B:318:LEU:HD12	1.60	0.82
1:A:10:TYR:CZ	1:A:84:LYS:HD3	2.16	0.81
1:A:482:THR:HG22	1:A:484:ASP:H	1.47	0.80
2:B:1099:ILE:HD12	2:B:1099:ILE:H	1.44	0.80
1:A:988:GLN:HG3	1:A:994:LYS:HB3	1.64	0.80
2:B:426:THR:OG1	2:B:427:VAL:CG2	2.30	0.80
1:A:525:ILE:HD12	1:A:525:ILE:H	1.48	0.79
2:B:336:LEU:HB3	2:B:347:MET:HE1	1.63	0.79
1:A:149:LEU:HD12	1:A:183:THR:CG2	2.11	0.79
3:C:18:LEU:HD22	3:C:20:LEU:HD12	1.64	0.79
1:A:80:THR:HG22	1:A:82:ASN:H	1.46	0.79
1:A:487:ILE:HG23	1:A:704:GLU:OE2	1.85	0.77
3:C:86:GLN:HB3	3:C:87:PRO:HD2	1.68	0.76
2:B:602:SER:HB2	2:B:608:GLU:OE1	1.85	0.75
1:A:52:THR:HG23	1:A:528:LEU:HD11	1.66	0.75
2:B:426:THR:HG21	2:B:428:ASP:OD2	1.86	0.75
2:B:336:LEU:HB3	2:B:347:MET:CE	2.18	0.74
1:A:1137:LYS:HG2	1:A:1138:GLY:H	1.53	0.74
1:A:421:GLY:N	1:A:424:GLU:HG3	2.02	0.74
2:B:1170:LYS:HE3	2:B:1173:ASN:HB3	1.69	0.73
2:B:1031:SER:O	2:B:1035:GLU:HG2	1.89	0.73
2:B:426:THR:OG1	2:B:427:VAL:HG22	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:251:GLU:HG2	2:B:357:HIS:HE1	1.54	0.72
3:C:18:LEU:HD23	3:C:19:ARG:N	2.05	0.72
1:A:149:LEU:HD12	1:A:183:THR:HG23	1.71	0.72
2:B:243:VAL:HG13	2:B:259:GLN:HB2	1.72	0.70
3:C:49:SER:HG	3:C:69:THR:HG1	1.33	0.70
1:A:1278:LEU:HD23	1:A:1281:SER:OG	1.92	0.70
1:A:1275:SER:HA	2:B:827:LEU:CD2	2.13	0.69
1:A:870:ILE:O	1:A:874:ILE:HG13	1.93	0.69
1:A:482:THR:HG22	1:A:484:ASP:N	2.07	0.69
2:B:715:ILE:HG22	2:B:720:LEU:HD22	1.76	0.68
2:B:243:VAL:CG1	2:B:259:GLN:HB2	2.22	0.68
2:B:1123:TYR:CE1	2:B:1138:LYS:HB2	2.28	0.68
2:B:834:TRP:HB2	2:B:835:PRO:HD2	1.74	0.68
2:B:96:GLU:HG2	2:B:459:ILE:HD13	1.75	0.67
2:B:426:THR:OG1	2:B:427:VAL:HG23	1.94	0.67
2:B:528:ASP:HB2	2:B:529:LYS:CA	2.25	0.66
2:B:528:ASP:HB2	2:B:529:LYS:HA	1.76	0.66
2:B:310:ASN:OD1	2:B:310:ASN:N	2.29	0.65
1:A:1228:ASP:HB3	1:A:1232:THR:H	1.60	0.65
2:B:295:PHE:CD1	2:B:328:ILE:HD11	2.32	0.65
2:B:340:SER:HB2	2:B:347:MET:CE	2.26	0.65
1:A:1236:LYS:NZ	1:A:1280:CYS:SG	2.70	0.64
3:C:1:GLN:HG3	3:C:2:VAL:H	1.62	0.64
2:B:493:ILE:HD11	2:B:499:TRP:CZ2	2.33	0.64
2:B:302:TYR:OH	2:B:321:LYS:HB2	1.98	0.64
1:A:149:LEU:CD1	1:A:183:THR:HG21	2.27	0.64
1:A:1170:LYS:H	1:A:1170:LYS:HD3	1.60	0.64
2:B:340:SER:HB2	2:B:347:MET:HE2	1.79	0.64
1:A:1170:LYS:N	1:A:1170:LYS:HD3	2.13	0.64
2:B:1132:LEU:HD21	2:B:1165:ILE:HD13	1.80	0.64
1:A:453:LEU:H	1:A:453:LEU:HD23	1.62	0.64
1:A:149:LEU:HD12	1:A:183:THR:HG21	1.79	0.63
1:A:681:LEU:HD23	1:A:694:THR:HG22	1.81	0.63
1:A:1:MET:N	1:A:2:PRO:HD3	2.14	0.63
2:B:409:ASN:OD1	2:B:409:ASN:N	2.30	0.63
1:A:611:VAL:HG12	1:A:1024:LEU:HD11	1.81	0.63
1:A:117:PHE:HA	1:A:317:PHE:CE1	2.34	0.62
1:A:80:THR:HG22	1:A:82:ASN:N	2.12	0.62
1:A:1271:ILE:HG13	2:B:854:ALA:HB1	1.81	0.62
2:B:251:GLU:HG2	2:B:357:HIS:CE1	2.34	0.62
1:A:1170:LYS:O	1:A:1170:LYS:HG2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:361:LYS:NZ	6:B:1196:SO4:O4	2.30	0.61
3:C:24:ALA:HB3	3:C:76:ASN:ND2	2.15	0.61
2:B:960:ASN:OD1	2:B:962:SER:HB3	2.00	0.61
1:A:948:ARG:NH1	8:A:1473:HOH:O	2.32	0.61
1:A:10:TYR:CZ	1:A:84:LYS:CD	2.84	0.61
1:A:135:ILE:HD13	1:A:149:LEU:CD2	2.30	0.61
1:A:52:THR:CG2	1:A:528:LEU:HD11	2.31	0.61
1:A:482:THR:CG2	1:A:484:ASP:H	2.14	0.60
1:A:53:ASN:OD1	1:A:55:GLU:HB2	2.01	0.60
2:B:567:ARG:NH2	8:B:1301:HOH:O	2.33	0.60
1:A:67:GLN:O	1:A:460:TRP:HZ2	1.85	0.60
2:B:1142:LYS:HD3	2:B:1160:TYR:CE1	2.36	0.60
3:C:97:THR:HG21	3:C:108:GLY:CA	2.31	0.60
1:A:1092:SER:O	1:A:1098:LYS:NZ	2.34	0.60
1:A:184:GLN:OE1	1:A:231:ARG:HD3	2.01	0.60
2:B:639:LEU:HD12	2:B:783:PHE:CE1	2.37	0.59
1:A:963:THR:HB	1:A:1058:ASP:HB3	1.84	0.59
1:A:706:TRP:CD2	1:A:808:LEU:HD13	2.37	0.59
1:A:149:LEU:CD1	1:A:183:THR:CG2	2.80	0.59
2:B:1125:ASP:OD1	2:B:1126:ILE:N	2.36	0.59
1:A:568:LEU:HD12	1:A:594:VAL:HG11	1.85	0.59
3:C:84:ASN:ND2	3:C:84:ASN:O	2.29	0.59
2:B:674:THR:HG22	2:B:675:GLN:HG3	1.85	0.59
2:B:79:LEU:O	2:B:83:ILE:HG13	2.03	0.59
1:A:388:ASP:HB3	1:A:391:ASN:O	2.03	0.59
2:B:924:LYS:HG3	2:B:928:CYS:O	2.03	0.59
2:B:313:GLY:H	2:B:317:LYS:HZ2	1.51	0.59
2:B:639:LEU:HD12	2:B:783:PHE:HE1	1.67	0.58
1:A:1128:VAL:HG11	1:A:1191:TYR:CE2	2.38	0.58
2:B:1082:ASN:ND2	2:B:1084:THR:O	2.36	0.58
2:B:169:ILE:CD1	2:B:199:THR:HB	2.33	0.58
2:B:1172:GLU:H	2:B:1174:HIS:H	1.50	0.58
3:C:30:GLY:HA2	3:C:71:ARG:CZ	2.34	0.58
1:A:1277:THR:HG21	2:B:341:LYS:HE3	1.85	0.58
1:A:1117:TYR:CD1	1:A:1277:THR:OG1	2.55	0.58
3:C:29:LEU:HD11	3:C:78:LEU:HB2	1.85	0.57
3:C:99:TYR:CD1	3:C:100:ARG:HB2	2.39	0.57
2:B:110:TYR:O	2:B:150:TYR:HD1	1.87	0.57
1:A:1228:ASP:OD2	1:A:1231:ILE:HB	2.05	0.56
1:A:687:ASN:ND2	1:A:690:LEU:HD22	2.19	0.56
2:B:281:TYR:OH	2:B:440:ARG:O	2.23	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:238:PRO:HG3	2:B:287:TYR:CZ	2.40	0.56
1:A:985:TRP:CD2	1:A:1019:ILE:HG21	2.40	0.56
2:B:169:ILE:HD13	2:B:199:THR:HB	1.88	0.56
1:A:1277:THR:O	1:A:1277:THR:HG23	2.05	0.56
1:A:67:GLN:HG2	1:A:425:PHE:CE1	2.41	0.56
1:A:869:TYR:HD1	1:A:870:ILE:HD12	1.72	0.55
2:B:283:PHE:HA	2:B:434:TYR:OH	2.06	0.55
1:A:80:THR:HB	1:A:83:GLU:HG3	1.88	0.55
3:C:97:THR:HG21	3:C:108:GLY:HA2	1.89	0.55
1:A:648:TYR:O	1:A:651:ASP:HB2	2.07	0.55
1:A:49:ASP:OD1	1:A:52:THR:HB	2.06	0.55
1:A:553:TYR:CZ	1:A:572:VAL:HG21	2.42	0.54
1:A:35:ALA:HB2	1:A:45:ILE:HG12	1.88	0.54
1:A:1271:ILE:HG22	1:A:1271:ILE:O	2.07	0.54
1:A:322:LEU:HD12	1:A:341:LEU:HB2	1.89	0.54
2:B:1137:ASN:OD1	2:B:1139:ASN:HB2	2.07	0.54
3:C:13:GLN:HB3	3:C:14:PRO:HD2	1.89	0.54
2:B:112:TYR:CE1	2:B:149:GLY:HA3	2.42	0.54
3:C:18:LEU:HD22	3:C:20:LEU:CD1	2.36	0.54
1:A:61:PRO:HB3	1:A:419:PHE:CZ	2.43	0.54
1:A:42:ILE:HD13	1:A:151:LEU:HB3	1.90	0.54
2:B:762:TYR:HB3	2:B:763:PRO:HD3	1.89	0.53
3:C:6:GLU:HA	3:C:21:SER:O	2.07	0.53
1:A:108:LEU:O	1:A:112:VAL:HG23	2.08	0.53
1:A:565:ARG:HB2	1:A:749:TYR:CZ	2.43	0.53
1:A:1228:ASP:HB3	1:A:1232:THR:N	2.23	0.53
2:B:306:ILE:HG23	2:B:312:ILE:CD1	2.39	0.53
2:B:227:ILE:HD11	2:B:339:PHE:CE2	2.43	0.53
2:B:513:LYS:HD2	2:B:514:PHE:H	1.74	0.53
3:C:24:ALA:HB3	3:C:76:ASN:HD21	1.74	0.53
2:B:513:LYS:HD2	2:B:514:PHE:N	2.23	0.53
1:A:1122:TYR:OH	1:A:1137:LYS:HD2	2.09	0.52
1:A:1241:ASP:HB3	1:A:1247:ILE:HD11	1.91	0.52
2:B:329:ASN:O	2:B:333:GLU:HG2	2.10	0.52
1:A:67:GLN:O	1:A:460:TRP:CZ2	2.62	0.52
1:A:1137:LYS:HG2	1:A:1138:GLY:N	2.21	0.52
2:B:715:ILE:CG2	2:B:720:LEU:HD22	2.38	0.52
1:A:1003:GLN:HA	1:A:1011:ILE:HD11	1.92	0.52
2:B:1043:ARG:HB3	2:B:1047:GLU:HA	1.92	0.52
1:A:115:ILE:HD11	1:A:512:PRO:HB2	1.92	0.52
1:A:207:LEU:HD11	1:A:1076:ASP:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:40:TRP:CD1	2:B:167:SER:HB2	2.45	0.52
3:C:37:VAL:HG23	3:C:46:GLU:O	2.10	0.52
1:A:227:HIS:CE1	1:A:262:GLU:OE2	2.63	0.52
1:A:1228:ASP:OD1	1:A:1234:LYS:HD2	2.09	0.52
2:B:341:LYS:HA	2:B:824:GLU:HG2	1.90	0.52
1:A:1209:LEU:CD1	1:A:1217:LEU:HD13	2.40	0.52
1:A:52:THR:HG22	1:A:53:ASN:N	2.25	0.52
2:B:266:LEU:O	2:B:335:ASN:HB3	2.10	0.52
2:B:887:ILE:HG23	2:B:1005:LEU:HB3	1.92	0.52
2:B:112:TYR:HE1	2:B:149:GLY:HA3	1.75	0.51
2:B:1058:LEU:HD23	2:B:1068:ILE:HG13	1.93	0.51
1:A:1278:LEU:HG	1:A:1280:CYS:SG	2.51	0.51
1:A:172:VAL:HG23	1:A:173:LEU:N	2.26	0.51
1:A:644:GLY:O	1:A:645:ASN:ND2	2.43	0.51
1:A:888:LEU:HB3	1:A:900:ILE:HD11	1.91	0.51
2:B:864:SER:HB2	2:B:867:ILE:HG13	1.93	0.51
1:A:699:LEU:HD22	1:A:844:THR:HG21	1.93	0.51
2:B:426:THR:CG2	2:B:428:ASP:OD2	2.58	0.51
1:A:671:ILE:O	1:A:671:ILE:HG23	2.11	0.51
1:A:864:SER:HA	1:A:867:THR:HB	1.93	0.51
3:C:67:PHE:N	3:C:67:PHE:CD1	2.77	0.51
2:B:336:LEU:O	2:B:347:MET:HE3	2.10	0.50
1:A:729:ARG:HD2	1:A:785:ASN:OD1	2.11	0.50
2:B:227:ILE:HG22	2:B:267:VAL:HG22	1.92	0.50
3:C:10:GLY:HA2	3:C:18:LEU:CD1	2.41	0.50
2:B:1:MET:HG2	2:B:2:ASN:H	1.75	0.50
1:A:1031:ILE:HG12	1:A:1036:ILE:HG13	1.92	0.50
1:A:1096:ILE:O	1:A:1098:LYS:HE3	2.11	0.50
1:A:913:GLN:HG2	1:A:1070:LYS:HD3	1.94	0.50
1:A:1238:ASN:HB2	1:A:1249:PHE:CE1	2.46	0.50
1:A:706:TRP:CE3	1:A:808:LEU:HD13	2.46	0.50
2:B:227:ILE:O	2:B:230:LEU:HB2	2.12	0.50
2:B:227:ILE:HD11	2:B:339:PHE:HE2	1.77	0.50
2:B:1172:GLU:H	2:B:1174:HIS:N	2.10	0.50
1:A:172:VAL:HG23	1:A:173:LEU:H	1.77	0.50
1:A:623:THR:HG23	6:A:1303:SO4:O1	2.12	0.49
2:B:615:ILE:HG13	2:B:615:ILE:O	2.12	0.49
1:A:177:ARG:HD3	1:A:238:ASN:HA	1.95	0.49
2:B:259:GLN:OE1	2:B:426:THR:O	2.31	0.49
1:A:45:ILE:HB	1:A:154:ILE:HG22	1.94	0.49
3:C:92:VAL:HG13	3:C:94:TYR:HE2	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:18:LEU:HD23	3:C:19:ARG:H	1.74	0.49
1:A:948:ARG:HB3	1:A:1068:TRP:HB2	1.95	0.49
2:B:874:ASN:HB2	2:B:1036:VAL:HG13	1.94	0.49
1:A:688:LYS:HD3	1:A:833:GLN:NE2	2.27	0.49
2:B:1115:SER:HB2	2:B:1123:TYR:CD2	2.48	0.48
2:B:374:ILE:HD12	2:B:375:ASN:N	2.27	0.48
2:B:326:ILE:HG23	2:B:327:ASN:N	2.28	0.48
2:B:528:ASP:CB	2:B:530:SER:N	2.61	0.48
1:A:634:ILE:N	1:A:634:ILE:HD12	2.27	0.48
2:B:633:GLU:HA	2:B:633:GLU:OE1	2.12	0.48
2:B:212:ASP:HB3	2:B:213:GLU:OE2	2.12	0.48
2:B:407:ASN:OD1	2:B:411:VAL:HG12	2.13	0.48
2:B:243:VAL:HG11	2:B:259:GLN:CD	2.33	0.48
1:A:23:LYS:HD2	3:C:33:TYR:OH	2.14	0.48
1:A:1186:VAL:HB	1:A:1191:TYR:CE1	2.49	0.48
1:A:532:PRO:O	1:A:534:ILE:HD12	2.14	0.47
2:B:1123:TYR:HD1	2:B:1138:LYS:HA	1.80	0.47
1:A:1164:LYS:NZ	1:A:1170:LYS:HA	2.29	0.47
2:B:316:ILE:HG22	2:B:317:LYS:N	2.29	0.47
2:B:320:LEU:HD12	2:B:320:LEU:HA	1.74	0.47
1:A:179:GLY:CA	1:A:234:GLY:HA3	2.44	0.47
2:B:1172:GLU:N	2:B:1173:ASN:HA	2.30	0.47
1:A:565:ARG:HB2	1:A:749:TYR:CE2	2.50	0.47
2:B:914:ASP:OD2	2:B:916:ILE:HG23	2.14	0.47
1:A:1014:TRP:CH2	1:A:1068:TRP:HB3	2.49	0.47
2:B:944:ILE:HB	2:B:950:GLU:HG3	1.95	0.47
1:A:858:ASP:OD1	1:A:858:ASP:N	2.47	0.47
2:B:1115:SER:HB2	2:B:1123:TYR:CE2	2.50	0.47
3:C:63:ALA:HA	3:C:66:ARG:HE	1.78	0.47
2:B:426:THR:CB	2:B:427:VAL:CA	2.69	0.47
1:A:525:ILE:N	1:A:525:ILE:HD12	2.23	0.47
2:B:528:ASP:CB	2:B:529:LYS:HA	2.40	0.47
2:B:272:ASP:N	2:B:273:PRO:HD2	2.29	0.47
2:B:966:ILE:HD12	2:B:966:ILE:N	2.30	0.47
2:B:902:PHE:HE1	2:B:904:ILE:HD11	1.80	0.46
1:A:590:TYR:O	1:A:594:VAL:HG12	2.15	0.46
2:B:238:PRO:HD2	2:B:263:VAL:CG2	2.45	0.46
3:C:90:THR:HG23	3:C:116:VAL:O	2.15	0.46
2:B:1166:CYS:HB3	2:B:1188:TYR:CD1	2.51	0.46
2:B:1039:ASN:O	2:B:1040:SER:CB	2.63	0.46
3:C:27:PHE:CE1	3:C:29:LEU:HD23	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:30:GLY:HA2	3:C:71:ARG:NH1	2.31	0.46
1:A:230:HIS:CE1	1:A:264:ARG:HD3	2.51	0.46
2:B:892:ASP:O	2:B:895:GLU:HG2	2.15	0.46
1:A:1125:VAL:HG22	1:A:1134:MET:HG2	1.98	0.46
1:A:97:ARG:HA	1:A:386:ILE:HG23	1.97	0.46
2:B:569:TYR:CE2	2:B:734:ILE:HG23	2.51	0.46
3:C:22:CYS:O	3:C:77:THR:HG23	2.16	0.46
2:B:1129:ASP:OD1	2:B:1129:ASP:N	2.49	0.46
2:B:92:THR:HB	2:B:343:PHE:CZ	2.51	0.46
1:A:80:THR:HG23	6:A:1306:SO4:O2	2.16	0.46
1:A:257:GLU:HG3	1:A:534:ILE:HG21	1.98	0.46
2:B:314:ASN:O	2:B:318:LEU:HB2	2.16	0.46
1:A:49:ASP:HB2	1:A:154:ILE:HD12	1.98	0.46
2:B:761:ILE:HD12	2:B:761:ILE:N	2.30	0.46
2:B:774:ASN:O	2:B:778:ILE:HG12	2.16	0.46
1:A:334:ASP:HB3	1:A:337:LYS:HB2	1.98	0.46
1:A:504:LEU:O	2:B:1170:LYS:HB2	2.16	0.45
2:B:1114:ILE:HB	2:B:1124:ILE:HD12	1.98	0.45
2:B:636:ASN:HA	2:B:639:LEU:CD1	2.47	0.45
1:A:26:ASN:HD21	1:A:54:PRO:HD3	1.81	0.45
2:B:587:LYS:HE2	2:B:587:LYS:HB3	1.67	0.45
1:A:22:ILE:HD11	1:A:45:ILE:HD11	1.99	0.45
2:B:155:TYR:HA	2:B:165:TYR:O	2.16	0.45
2:B:435:LYS:HA	2:B:628:ILE:HG23	1.98	0.45
1:A:838:LYS:HD3	8:A:1448:HOH:O	2.16	0.45
2:B:842:ALA:HA	2:B:850:VAL:O	2.16	0.45
1:A:773:LEU:O	1:A:777:ILE:HG13	2.17	0.45
2:B:184:VAL:CG2	2:B:200:GLU:HB2	2.47	0.45
1:A:930:ASN:O	1:A:933:VAL:HG13	2.16	0.45
3:C:2:VAL:HG13	3:C:27:PHE:CE2	2.51	0.45
2:B:306:ILE:HG23	2:B:312:ILE:HD12	1.99	0.45
2:B:17:ASN:HD22	2:B:34:LYS:HE3	1.81	0.45
7:B:1204:MES:H81	7:B:1204:MES:H52	1.71	0.45
2:B:1080:THR:OG1	2:B:1081:ILE:N	2.49	0.45
1:A:140:PRO:O	3:C:97:THR:HG23	2.16	0.45
2:B:427:VAL:HA	2:B:428:ASP:HA	1.68	0.44
3:C:10:GLY:H	3:C:18:LEU:HD11	1.82	0.44
2:B:261:ASN:OD1	2:B:263:VAL:HG23	2.17	0.44
1:A:118:TRP:HB3	1:A:128:LYS:O	2.17	0.44
1:A:125:THR:O	1:A:301:LYS:N	2.48	0.44
1:A:421:GLY:H	1:A:424:GLU:CG	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1209:LEU:HD11	1:A:1217:LEU:HD13	1.99	0.44
2:B:406:LEU:HD11	2:B:494:ASN:HB3	1.99	0.44
2:B:426:THR:HG1	2:B:427:VAL:HG22	1.82	0.44
1:A:1097:LEU:HD12	1:A:1225:SER:HB3	2.00	0.44
1:A:1193:LEU:HD11	1:A:1206:LEU:HD13	2.00	0.44
2:B:1156:LEU:HD21	2:B:1192:PHE:CD2	2.53	0.44
1:A:149:LEU:HD11	1:A:183:THR:HG21	2.00	0.44
2:B:1167:LEU:HD12	2:B:1167:LEU:HA	1.81	0.44
1:A:1155:TYR:CD2	1:A:1293:GLU:HG3	2.53	0.44
2:B:293:LYS:HB2	2:B:293:LYS:NZ	2.32	0.44
2:B:841:TYR:CZ	2:B:852:GLY:HA3	2.53	0.44
2:B:105:ALA:HB3	2:B:233:LEU:HD13	2.00	0.44
1:A:1122:TYR:CD1	1:A:1157:GLY:HA2	2.53	0.43
2:B:1123:TYR:CD1	2:B:1138:LYS:HA	2.53	0.43
3:C:27:PHE:HE1	3:C:29:LEU:HD23	1.83	0.43
2:B:154:ASN:HB2	2:B:167:SER:O	2.18	0.43
2:B:34:LYS:HB2	2:B:40:TRP:CZ3	2.53	0.43
2:B:46:TYR:OH	2:B:200:GLU:OE2	2.23	0.43
1:A:43:TRP:CD1	1:A:149:LEU:HD22	2.54	0.43
3:C:99:TYR:CE1	3:C:100:ARG:HB2	2.54	0.43
2:B:967:THR:HB	2:B:980:PHE:HB2	2.00	0.43
2:B:911:LEU:HG	2:B:911:LEU:H	1.68	0.43
1:A:501:GLN:HG3	2:B:1063:PHE:CE1	2.53	0.43
2:B:184:VAL:HG22	2:B:200:GLU:HB2	2.01	0.43
1:A:1202:VAL:HG13	1:A:1203:GLU:OE2	2.18	0.43
2:B:859:THR:HA	2:B:890:SER:O	2.18	0.43
3:C:14:PRO:HA	3:C:15:GLY:HA3	1.88	0.43
2:B:762:TYR:O	2:B:765:PHE:HB3	2.19	0.43
1:A:128:LYS:HG2	1:A:304:VAL:HB	1.99	0.43
1:A:711:LYS:HB3	1:A:711:LYS:HE2	1.80	0.43
2:B:1108:LYS:O	2:B:1109:LEU:HB2	2.19	0.43
3:C:72:ASP:OD2	3:C:75:LYS:HG3	2.18	0.43
2:B:426:THR:HB	2:B:428:ASP:HA	2.00	0.43
3:C:10:GLY:HA2	3:C:18:LEU:HD11	2.00	0.43
1:A:1134:MET:HB2	1:A:1134:MET:HE3	1.81	0.43
1:A:178:ASN:OD1	1:A:180:TYR:HB2	2.19	0.43
1:A:135:ILE:HD13	1:A:149:LEU:HD23	1.99	0.43
1:A:135:ILE:HD13	1:A:149:LEU:HD21	2.01	0.43
2:B:1172:GLU:HB2	2:B:1173:ASN:C	2.39	0.43
2:B:567:ARG:NH1	2:B:973:LEU:O	2.52	0.43
2:B:249:ARG:NH1	2:B:256:GLU:OE2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1271:ILE:HG13	2:B:854:ALA:CB	2.49	0.43
1:A:907:ASP:HB3	1:A:910:ASP:O	2.19	0.43
2:B:1100:ASN:HA	2:B:1101:PRO:HD2	1.85	0.43
1:A:600:ALA:HA	1:A:754:TYR:CZ	2.54	0.43
1:A:195:GLY:HA2	1:A:213:PHE:O	2.19	0.43
3:C:20:LEU:HD13	3:C:82:MET:SD	2.58	0.43
3:C:1:GLN:HB3	3:C:3:GLN:OE1	2.19	0.43
1:A:67:GLN:HA	1:A:425:PHE:CZ	2.54	0.43
3:C:63:ALA:HB1	3:C:67:PHE:CG	2.54	0.43
1:A:1030:TYR:O	1:A:1031:ILE:HD13	2.19	0.43
1:A:1023:ARG:NH1	1:A:1023:ARG:HB3	2.34	0.43
1:A:881:LEU:HD11	1:A:898:ILE:HG12	2.01	0.43
1:A:1293:GLU:O	1:A:1295:PRO:HD3	2.19	0.42
1:A:481:ILE:O	1:A:682:VAL:HG12	2.18	0.42
2:B:1173:ASN:O	2:B:1174:HIS:CD2	2.72	0.42
2:B:838:LEU:HD13	2:B:859:THR:HG21	2.01	0.42
1:A:164:GLU:OE2	1:A:166:LYS:HE2	2.19	0.42
2:B:1071:VAL:CG1	2:B:1176:TRP:HZ3	2.33	0.42
2:B:336:LEU:C	2:B:347:MET:HE3	2.40	0.42
2:B:554:THR:OG1	2:B:557:LYS:HG3	2.19	0.42
1:A:1188:ASN:ND2	1:A:1188:ASN:O	2.42	0.42
1:A:1135:TYR:HB2	1:A:1259:ALA:O	2.18	0.42
1:A:1245:ASN:ND2	2:B:846:PRO:HA	2.34	0.42
1:A:135:ILE:CD1	1:A:149:LEU:HD23	2.50	0.42
2:B:1123:TYR:HE1	2:B:1138:LYS:HB2	1.77	0.42
1:A:1:MET:H1	1:A:2:PRO:HD3	1.83	0.42
2:B:313:GLY:H	2:B:317:LYS:NZ	2.17	0.42
1:A:788:LEU:HD23	1:A:788:LEU:HA	1.82	0.42
2:B:709:LEU:HD12	2:B:709:LEU:HA	1.86	0.42
2:B:871:TYR:CE2	2:B:1049:ARG:HD3	2.55	0.42
1:A:1228:ASP:H	1:A:1232:THR:HA	1.85	0.42
2:B:636:ASN:HA	2:B:639:LEU:HD13	2.02	0.42
2:B:674:THR:HG22	2:B:675:GLN:N	2.35	0.42
1:A:985:TRP:CE2	1:A:1019:ILE:HG21	2.55	0.42
2:B:700:ARG:NH1	8:B:1275:HOH:O	2.52	0.42
1:A:1179:ARG:NH2	1:A:1240:GLN:OE1	2.50	0.42
2:B:1031:SER:O	2:B:1035:GLU:CG	2.64	0.42
2:B:20:VAL:HG13	2:B:29:VAL:HG13	2.01	0.42
2:B:1052:TYR:CE1	2:B:1098:SER:HB2	2.54	0.42
2:B:1084:THR:HB	2:B:1087:LEU:HB2	2.01	0.42
1:A:1277:THR:HG21	2:B:341:LYS:CE	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:471:PHE:CE2	1:A:720:LYS:HE3	2.54	0.42
1:A:1159:LYS:HB2	1:A:1185:VAL:HB	2.01	0.42
2:B:638:MET:HA	2:B:641:LEU:HD13	2.01	0.41
2:B:1144:MET:HB2	2:B:1144:MET:HE2	1.91	0.41
3:C:13:GLN:NE2	8:C:241:HOH:O	2.53	0.41
2:B:211:TYR:O	2:B:212:ASP:HB2	2.20	0.41
2:B:330:ASP:HB3	2:B:453:SER:HB3	2.02	0.41
2:B:185:PHE:HB2	2:B:197:THR:HG21	2.01	0.41
2:B:536:LEU:O	2:B:540:MET:HE2	2.21	0.41
2:B:426:THR:CB	2:B:427:VAL:CG2	2.97	0.41
1:A:10:TYR:CE1	1:A:84:LYS:HD3	2.53	0.41
1:A:10:TYR:OH	1:A:84:LYS:HD3	2.20	0.41
2:B:368:TYR:HA	2:B:369:PRO:HA	1.62	0.41
2:B:906:PHE:HB3	2:B:1017:LEU:HA	2.02	0.41
1:A:688:LYS:O	1:A:692:VAL:HG23	2.21	0.41
2:B:1122:LYS:HB3	2:B:1134:LEU:HB3	2.00	0.41
2:B:40:TRP:HD1	2:B:167:SER:HB2	1.84	0.41
2:B:1070:GLU:CD	2:B:1077:ILE:HD11	2.40	0.41
2:B:624:LEU:HA	2:B:624:LEU:HD12	1.76	0.41
1:A:964:ILE:HD13	1:A:978:LEU:HG	2.03	0.41
1:A:242:VAL:HG13	1:A:258:VAL:O	2.20	0.41
1:A:97:ARG:HD2	1:A:386:ILE:O	2.20	0.41
1:A:1023:ARG:NH2	1:A:1045:GLY:O	2.54	0.41
1:A:1111:TYR:CG	1:A:1284:PHE:HB3	2.56	0.41
2:B:648:GLU:HG2	2:B:648:GLU:O	2.19	0.41
2:B:177:ASN:ND2	2:B:180:GLU:HB2	2.36	0.41
3:C:4:LEU:O	3:C:111:GLY:HA2	2.20	0.41
2:B:1127:SER:OG	2:B:1128:GLU:N	2.54	0.41
2:B:579:ILE:O	2:B:585:ILE:HA	2.21	0.40
1:A:525:ILE:CD1	1:A:525:ILE:H	2.24	0.40
1:A:1:MET:N	1:A:2:PRO:CD	2.83	0.40
1:A:1:MET:HG2	1:A:1:MET:O	2.21	0.40
2:B:1152:ILE:CG2	2:B:1155:CYS:HB2	2.52	0.40
1:A:1120:ASN:O	1:A:1140:ARG:HG3	2.22	0.40
2:B:754:ILE:O	2:B:758:GLU:HG3	2.21	0.40
2:B:639:LEU:CD1	2:B:783:PHE:HE1	2.32	0.40
3:C:46:GLU:OE2	3:C:62:SER:HB2	2.19	0.40
1:A:634:ILE:HA	1:A:635:PRO:HD3	1.84	0.40
1:A:179:GLY:HA2	1:A:234:GLY:HA3	2.02	0.40
1:A:565:ARG:HD2	1:A:749:TYR:CE1	2.55	0.40
2:B:368:TYR:N	2:B:368:TYR:CD2	2.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1130:ASN:HB3	2:B:1179:CYS:O	2.21	0.40
1:A:1277:THR:O	1:A:1277:THR:CG2	2.67	0.40
2:B:511:ASN:HB3	2:B:513:LYS:O	2.21	0.40
1:A:266:PHE:CE1	1:A:366:PHE:HB3	2.56	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	1276/1296 (98%)	1242 (97%)	34 (3%)	0	100	100
2	B	1144/1196 (96%)	1105 (97%)	38 (3%)	1 (0%)	56	83
3	C	117/152 (77%)	112 (96%)	5 (4%)	0	100	100
All	All	2537/2644 (96%)	2459 (97%)	77 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	429	ASP

### 5.3.2 Protein sidechains [i](#)

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	1162/1176 (99%)	1125 (97%)	37 (3%)	46	77
2	B	1076/1114 (97%)	1008 (94%)	68 (6%)	22	48
3	C	94/123 (76%)	83 (88%)	11 (12%)	7	15
All	All	2332/2413 (97%)	2216 (95%)	116 (5%)	30	60

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	THR
1	A	55	GLU
1	A	58	ASP
1	A	97	ARG
1	A	240	ASN
1	A	482	THR
1	A	484	ASP
1	A	507	ASN
1	A	510	ASN
1	A	520	LEU
1	A	525	ILE
1	A	541	LYS
1	A	546	ASP
1	A	608	GLU
1	A	645	ASN
1	A	682	VAL
1	A	690	LEU
1	A	757	GLU
1	A	770	SER
1	A	806	LYS
1	A	819	LEU
1	A	858	ASP
1	A	871	LYS
1	A	933	VAL
1	A	948	ARG
1	A	958	LEU
1	A	1027	SER
1	A	1065	ARG
1	A	1087	LEU
1	A	1146	THR
1	A	1156	ARG
1	A	1169	ASN
1	A	1170	LYS
1	A	1188	ASN

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Mol	Chain	Res	Type
1	A	1226	LYS
1	A	1272	GLU
1	A	1274	SER
2	B	7	LEU
2	B	29	VAL
2	B	50	SER
2	B	186	TYR
2	B	230	LEU
2	B	250	SER
2	B	251	GLU
2	B	253	GLU
2	B	279	ASP
2	B	284	THR
2	B	310	ASN
2	B	320	LEU
2	B	326	ILE
2	B	329	ASN
2	B	359	TYR
2	B	406	LEU
2	B	409	ASN
2	B	469	ILE
2	B	487	ILE
2	B	492	GLU
2	B	493	ILE
2	B	504	LEU
2	B	509	THR
2	B	511	ASN
2	B	528	ASP
2	B	550	SER
2	B	553	ASP
2	B	554	THR
2	B	583	CYS
2	B	585	ILE
2	B	611	ASN
2	B	621	LYS
2	B	624	LEU
2	B	648	GLU
2	B	674	THR
2	B	703	GLN
2	B	709	LEU
2	B	712	ASN
2	B	716	SER

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Mol	Chain	Res	Type
2	B	720	LEU
2	B	724	ASN
2	B	745	VAL
2	B	796	LEU
2	B	797	GLN
2	B	800	ASN
2	B	830	LYS
2	B	834	TRP
2	B	865	LYS
2	B	913	LYS
2	B	916	ILE
2	B	924	LYS
2	B	927	ASN
2	B	978	LEU
2	B	993	GLU
2	B	1024	THR
2	B	1026	SER
2	B	1077	ILE
2	B	1083	ASN
2	B	1086	ASN
2	B	1089	LEU
2	B	1115	SER
2	B	1117	LEU
2	B	1128	GLU
2	B	1129	ASP
2	B	1136	ASP
2	B	1142	LYS
2	B	1144	MET
2	B	1167	LEU
3	C	0	SER
3	C	5	VAL
3	C	11	LEU
3	C	12	VAL
3	C	28	THR
3	C	29	LEU
3	C	49	SER
3	C	59	ASP
3	C	76	ASN
3	C	84	ASN
3	C	88	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	913	GLN
2	B	961	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 26 ligands modelled in this entry, 2 are monoatomic - leaving 24 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	$\# Z  > 2$
6	SO4	A	1299	-	4,4,4	0.19	0	6,6,6	0.11	0
6	SO4	A	1300	-	4,4,4	0.19	0	6,6,6	0.10	0
6	SO4	A	1301	-	4,4,4	0.17	0	6,6,6	0.19	0
6	SO4	A	1302	-	4,4,4	0.21	0	6,6,6	0.10	0
6	SO4	A	1303	-	4,4,4	0.22	0	6,6,6	0.19	0
6	SO4	A	1304	-	4,4,4	0.21	0	6,6,6	0.10	0
6	SO4	A	1305	-	4,4,4	0.20	0	6,6,6	0.08	0
6	SO4	A	1306	-	4,4,4	0.21	0	6,6,6	0.11	0
6	SO4	A	1307	-	4,4,4	0.21	0	6,6,6	0.11	0
6	SO4	A	1308	-	4,4,4	0.19	0	6,6,6	0.15	0
6	SO4	A	1309	-	4,4,4	0.20	0	6,6,6	0.06	0
6	SO4	A	1310	-	4,4,4	0.18	0	6,6,6	0.13	0
6	SO4	B	1195	-	4,4,4	0.19	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	SO4	B	1196	-	4,4,4	0.20	0	6,6,6	0.08	0
6	SO4	B	1197	-	4,4,4	0.22	0	6,6,6	0.13	0
6	SO4	B	1198	-	4,4,4	0.15	0	6,6,6	0.11	0
6	SO4	B	1199	-	4,4,4	0.19	0	6,6,6	0.14	0
6	SO4	B	1200	-	4,4,4	0.21	0	6,6,6	0.16	0
6	SO4	B	1201	-	4,4,4	0.21	0	6,6,6	0.15	0
7	MES	B	1202	-	11,12,12	1.13	1 (9%)	14,16,16	2.38	3 (21%)
7	MES	B	1203	-	11,12,12	1.21	1 (9%)	14,16,16	2.34	3 (21%)
7	MES	B	1204	-	11,12,12	1.18	1 (9%)	14,16,16	2.52	3 (21%)
7	MES	B	1205	-	11,12,12	1.18	1 (9%)	14,16,16	2.45	3 (21%)
6	SO4	C	144	-	4,4,4	0.20	0	6,6,6	0.08	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
6	SO4	A	1299	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1300	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1301	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1302	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1303	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1304	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1305	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1306	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1307	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1308	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1309	-	-	0/0/0/0	0/0/0/0
6	SO4	A	1310	-	-	0/0/0/0	0/0/0/0
6	SO4	B	1195	-	-	0/0/0/0	0/0/0/0
6	SO4	B	1196	-	-	0/0/0/0	0/0/0/0
6	SO4	B	1197	-	-	0/0/0/0	0/0/0/0
6	SO4	B	1198	-	-	0/0/0/0	0/0/0/0
6	SO4	B	1199	-	-	0/0/0/0	0/0/0/0
6	SO4	B	1200	-	-	0/0/0/0	0/0/0/0
6	SO4	B	1201	-	-	0/0/0/0	0/0/0/0
7	MES	B	1202	-	-	0/6/14/14	0/1/1/1
7	MES	B	1203	-	-	0/6/14/14	0/1/1/1
7	MES	B	1204	-	-	0/6/14/14	0/1/1/1
7	MES	B	1205	-	-	0/6/14/14	0/1/1/1
6	SO4	C	144	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	1205	MES	C5-N4	-3.10	1.38	1.46
7	B	1203	MES	C5-N4	-3.08	1.38	1.46
7	B	1204	MES	C5-N4	-3.05	1.38	1.46
7	B	1202	MES	C5-N4	-2.89	1.39	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1204	MES	O3S-S-O2S	-4.67	100.73	111.61
7	B	1202	MES	O3S-S-O2S	-4.54	101.04	111.61
7	B	1205	MES	O3S-S-O2S	-4.52	101.09	111.61
7	B	1203	MES	O3S-S-O2S	-4.22	101.78	111.61
7	B	1203	MES	O1S-S-C8	3.74	110.10	106.91
7	B	1202	MES	O1S-S-C8	4.42	110.68	106.91
7	B	1204	MES	O1S-S-C8	4.71	110.92	106.91
7	B	1205	MES	O1S-S-C8	4.91	111.09	106.91
7	B	1202	MES	O2S-S-C8	4.97	111.15	106.91
7	B	1205	MES	O2S-S-C8	5.22	111.36	106.91
7	B	1203	MES	O2S-S-C8	5.62	111.70	106.91
7	B	1204	MES	O2S-S-C8	5.69	111.76	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1303	SO4	1	0
6	A	1306	SO4	1	0
6	B	1196	SO4	1	0
7	B	1204	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	1280/1296 (98%)	-0.24	16 (1%) 79 79	32, 56, 101, 170	10 (0%)
2	B	1150/1196 (96%)	-0.12	24 (2%) 67 68	29, 54, 111, 166	14 (1%)
3	C	119/152 (78%)	1.09	27 (22%) 1 1	79, 115, 158, 187	1 (0%)
All	All	2549/2644 (96%)	-0.12	67 (2%) 59 59	29, 56, 117, 187	25 (0%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	9	GLY	5.8
3	C	14	PRO	5.8
1	A	450	LEU	4.7
3	C	93	TYR	4.3
3	C	118	VAL	4.2
3	C	10	GLY	4.1
1	A	1276	ARG	4.0
3	C	85	LEU	3.9
3	C	19	ARG	3.8
2	B	1086	ASN	3.7
2	B	112	TYR	3.6
3	C	83	SER	3.6
3	C	18	LEU	3.5
3	C	67	PHE	3.4
2	B	638	MET	3.4
2	B	635	PRO	3.4
3	C	20	LEU	3.3
2	B	149	GLY	3.3
2	B	150	TYR	3.2
3	C	87	PRO	3.1
3	C	61	ASP	3.1
2	B	634	ILE	3.1
1	A	1214	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	452	ASP	3.0
2	B	1085	ASN	2.9
1	A	1130	ILE	2.9
3	C	82	MET	2.9
3	C	38	ARG	2.9
3	C	13	GLN	2.9
2	B	312	ILE	2.9
3	C	16	GLY	2.9
2	B	412	SER	2.8
2	B	451	ASP	2.8
1	A	1	MET	2.8
1	A	1275	SER	2.8
2	B	113	ILE	2.7
2	B	637	ASP	2.7
2	B	441	ALA	2.7
1	A	1200	ALA	2.7
3	C	116	VAL	2.7
3	C	21	SER	2.7
2	B	306	ILE	2.6
1	A	451	ASN	2.6
3	C	51	ILE	2.6
2	B	409	ASN	2.6
3	C	79	TYR	2.5
2	B	1088	ASN	2.5
1	A	646	MET	2.4
3	C	69	THR	2.4
2	B	783	PHE	2.4
2	B	111	GLY	2.4
3	C	66	ARG	2.3
3	C	11	LEU	2.3
3	C	59	ASP	2.3
2	B	1077	ILE	2.2
2	B	301	ILE	2.2
3	C	114	THR	2.2
1	A	1197	ALA	2.2
2	B	641	LEU	2.1
1	A	491	GLU	2.1
2	B	411	VAL	2.1
1	A	453	LEU	2.1
1	A	485	THR	2.0
2	B	632	TYR	2.0
3	C	33	TYR	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	433	GLY	2.0
1	A	1294	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
7	MES	B	1205	12/12	0.85	0.28	12.64	105,116,150,150	0
6	SO4	B	1199	5/5	0.87	0.35	9.63	122,125,128,130	0
6	SO4	B	1196	5/5	0.88	0.27	6.40	157,159,161,162	0
6	SO4	A	1310	5/5	0.88	0.30	6.25	133,139,139,142	0
7	MES	B	1204	12/12	0.83	0.22	6.20	82,100,176,178	0
6	SO4	A	1307	5/5	0.83	0.36	5.74	172,173,174,176	0
6	SO4	A	1301	5/5	0.90	0.24	5.10	120,121,125,125	0
6	SO4	A	1303	5/5	0.85	0.23	4.05	132,136,137,140	0
6	SO4	B	1201	5/5	0.86	0.20	2.96	151,152,155,160	0
6	SO4	A	1304	5/5	0.90	0.16	2.47	136,137,138,139	0
6	SO4	A	1309	5/5	0.85	0.23	2.43	159,159,160,160	0
6	SO4	A	1302	5/5	0.89	0.26	0.86	137,137,138,139	0
7	MES	B	1202	12/12	0.84	0.21	0.80	88,96,189,190	0
6	SO4	B	1195	5/5	0.93	0.14	0.31	96,102,104,104	0
7	MES	B	1203	12/12	0.96	0.17	-0.26	74,79,111,112	0
6	SO4	A	1299	5/5	0.98	0.16	-0.26	97,99,101,103	0
5	CA	A	1298	1/1	0.97	0.05	-4.89	67,67,67,67	1
6	SO4	B	1198	5/5	0.95	0.15	-	86,94,105,105	0
4	ZN	A	1297	1/1	0.99	0.17	-	62,62,62,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	SO4	A	1306	5/5	0.78	0.26	-	157,160,162,162	0
6	SO4	A	1305	5/5	0.81	0.30	-	155,157,160,161	0
6	SO4	B	1200	5/5	0.90	0.15	-	112,117,119,121	0
6	SO4	A	1300	5/5	0.96	0.14	-	98,103,108,110	0
6	SO4	C	144	5/5	0.78	0.22	-	161,162,165,165	0
6	SO4	B	1197	5/5	0.93	0.22	-	129,131,132,135	0
6	SO4	A	1308	5/5	0.95	0.21	-	139,142,142,143	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.