



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 10:42 PM EDT

PDB ID : 3WU3  
Title : Reduced-form structure of E.coli Lon Proteolytic domain  
Authors : Nishii, W.; Kukimoto-Niino, M.; Terada, T.; Shirouzu, M.; Muramatsu, T.;  
Yokoyama, S.  
Deposited on : 2014-04-22  
Resolution : 1.82 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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.20.1  
EDS : 2.36.2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36.2

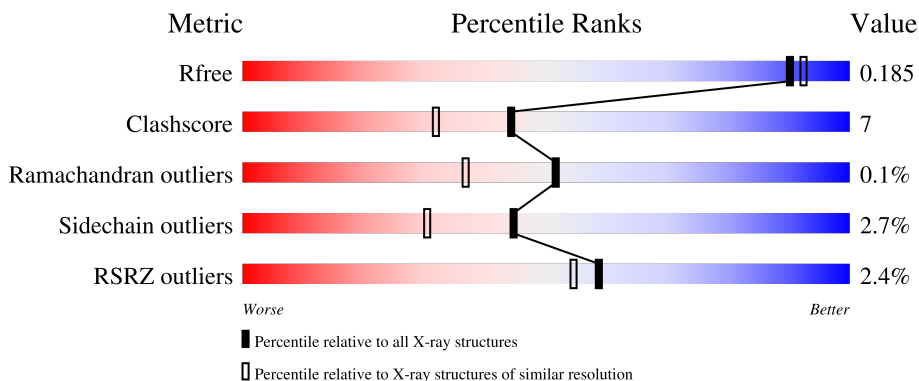
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.82 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.




Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	7484 (1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	200	 84% 10% 5%
1	B	200	 4% 76% 13% 10%
1	C	200	 73% 17% 10%
1	D	200	 3% 72% 17% 10%
1	E	200	 4% 74% 22% 5%

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Mol	Chain	Length	Quality of chain
1	F	200	 76% 14% 10%

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8431 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 Lon protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	190	1425	894	251	274	6	0	0	0
1	B	181	1349	851	234	258	6	0	0	0
1	C	181	1349	851	234	258	6	0	0	0
1	D	181	1349	851	234	258	6	0	0	0
1	E	190	1412	889	245	271	7	0	0	0
1	F	179	1333	841	232	254	6	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	679	ALA	SER	engineered mutation	UNP C9QQ79
B	679	ALA	SER	engineered mutation	UNP C9QQ79
C	679	ALA	SER	engineered mutation	UNP C9QQ79
D	679	ALA	SER	engineered mutation	UNP C9QQ79
E	679	ALA	SER	engineered mutation	UNP C9QQ79
F	679	ALA	SER	engineered mutation	UNP C9QQ79

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

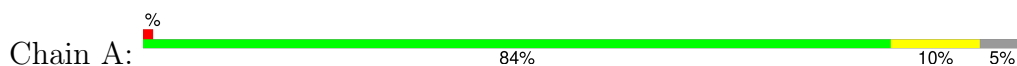
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	40	Total	O	0	0
			40	40		
3	B	26	Total	O	0	0
			26	26		
3	C	32	Total	O	0	0
			32	32		
3	D	26	Total	O	0	0
			26	26		
3	E	32	Total	O	0	0
			32	32		
3	F	33	Total	O	0	0
			33	33		

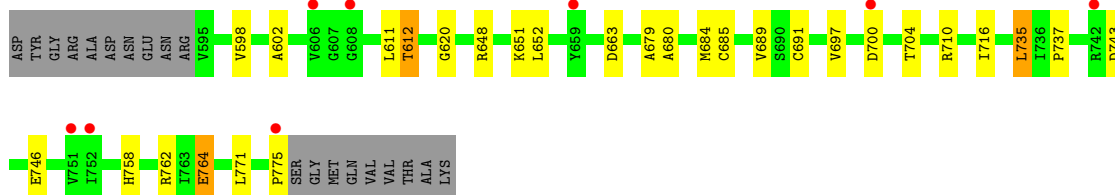
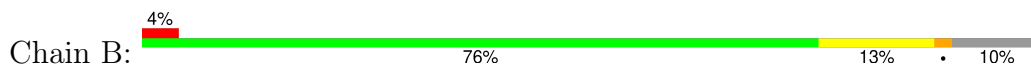
### 3 Residue-property plots [i](#)

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

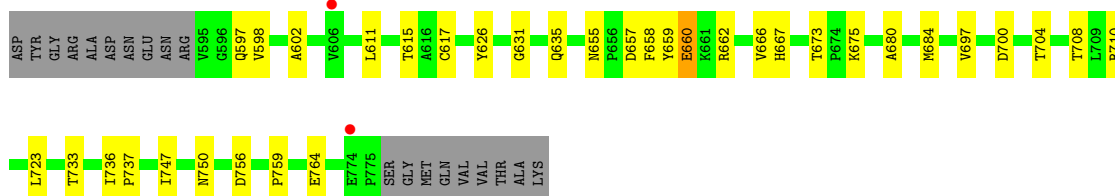
- Molecule 1: Lon protease



- Molecule 1: Lon protease



- Molecule 1: Lon protease

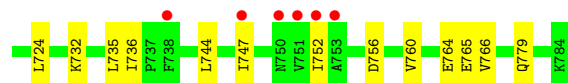


- Molecule 1: Lon protease

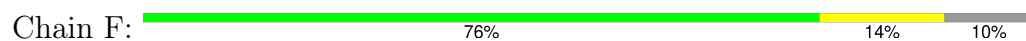




- Molecule 1: Lon protease



- Molecule 1: Lon protease



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.24Å 86.24Å 123.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.37 – 1.82 35.37 – 1.82	Depositor EDS
% Data completeness (in resolution range)	98.7 (35.37-1.82) 91.9 (35.37-1.82)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.64 (at 1.82Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.1_1168)	Depositor
R, $R_{free}$	0.160 , 0.196 0.159 , 0.185	Depositor DCC
$R_{free}$ test set	4536 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.8	Xtrriage
Anisotropy	0.773	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 54.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.22$	Xtrriage
Estimated twinning fraction	0.268 for -h,-k,l 0.105 for h,-h-k,-l 0.098 for -k,-h,-l	Xtrriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	8431	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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: 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	0.38	0/1446	0.54	0/1964
1	B	0.35	0/1369	0.53	0/1861
1	C	0.35	0/1369	0.53	0/1861
1	D	0.34	0/1369	0.55	0/1861
1	E	0.35	0/1432	0.53	0/1944
1	F	0.36	0/1352	0.53	0/1837
All	All	0.36	0/8337	0.54	0/11328

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 [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	1425	0	1458	16	0
1	B	1349	0	1393	17	0
1	C	1349	0	1393	20	0
1	D	1349	0	1393	20	0
1	E	1412	0	1461	25	0
1	F	1333	0	1380	21	0
2	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	40	0	0	0	0
3	B	26	0	0	0	0
3	C	32	0	0	3	0
3	D	26	0	0	0	0
3	E	32	0	0	2	0
3	F	33	0	0	1	0
All	All	8431	0	8478	110	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 7.

All (110) 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:E:658:PHE:HA	1:E:661:LYS:HE2	1.59	0.82
1:C:673:THR:O	1:C:675:LYS:NZ	2.14	0.79
1:F:723:LEU:HD21	1:F:734:VAL:HG11	1.71	0.71
1:C:750:ASN:OD1	3:C:932:HOH:O	2.11	0.68
1:F:706:GLU:OE2	3:F:922:HOH:O	2.12	0.66
1:D:628:GLY:O	1:D:675:LYS:NZ	2.24	0.66
1:F:698:ARG:HG2	1:F:772:GLN:HG2	1.78	0.65
1:C:602:ALA:HB3	1:C:611:LEU:HD12	1.78	0.64
1:A:586:TYR:CZ	1:B:775:PRO:HG2	2.32	0.64
1:E:700:ASP:HB2	1:E:732:LYS:HG2	1.80	0.64
1:C:598:VAL:HG22	1:C:697:VAL:HG11	1.79	0.63
1:A:617:CYS:SG	1:A:662:ARG:HD3	2.39	0.63
1:C:710:ARG:NH2	3:C:917:HOH:O	2.31	0.61
1:E:673:THR:O	1:E:675:LYS:NZ	2.29	0.60
1:A:601:LEU:HB2	1:A:704:THR:HB	1.83	0.59
1:B:771:LEU:HD12	1:B:775:PRO:HB3	1.82	0.59
1:E:619:PRO:HA	1:E:662:ARG:HD3	1.85	0.58
1:E:615:THR:HG22	1:E:666:VAL:HG22	1.86	0.57
1:B:598:VAL:HG22	1:B:697:VAL:HG21	1.86	0.57
1:D:615:THR:HG22	1:D:666:VAL:HG13	1.87	0.56
1:E:732:LYS:O	1:E:756:ASP:N	2.30	0.56
1:E:747:ILE:HB	1:E:752:ILE:HD11	1.87	0.55
1:D:685:CYS:O	1:D:689:VAL:HG23	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:631:GLY:O	1:E:635:GLN:HG3	2.07	0.54
1:A:586:TYR:O	1:B:648:ARG:NH1	2.38	0.54
1:D:742:ARG:NH1	1:D:743:ASP:OD1	2.41	0.54
1:E:655:ASN:ND2	1:E:657:ASP:HB2	2.24	0.53
1:C:615:THR:HG22	1:C:666:VAL:HG22	1.90	0.52
1:C:723:LEU:HD12	1:C:747:ILE:HD13	1.92	0.52
1:C:667:HIS:HB3	1:D:709:LEU:HG	1.91	0.52
1:C:597:GLN:OE1	1:D:710:ARG:HD3	2.10	0.52
1:D:603:TRP:HA	1:D:608:GLY:HA2	1.92	0.52
1:F:723:LEU:HD12	1:F:747:ILE:HD13	1.90	0.52
1:E:614:GLU:OE2	1:F:710:ARG:NH1	2.43	0.51
1:E:655:ASN:HD21	1:E:657:ASP:HB2	1.76	0.51
1:A:615:THR:HG22	1:A:666:VAL:HG22	1.92	0.50
1:C:631:GLY:O	1:C:635:GLN:HG3	2.11	0.50
1:C:704:THR:O	1:C:737:PRO:HD3	2.12	0.50
1:D:601:LEU:HB2	1:D:704:THR:HB	1.93	0.50
1:D:716:ILE:HD11	1:D:719:LEU:HD13	1.94	0.49
1:D:689:VAL:O	1:D:693:THR:OG1	2.23	0.49
1:E:640:ALA:O	1:E:644:VAL:HG23	2.12	0.49
1:B:679:ALA:HA	1:B:716:ILE:HG21	1.94	0.48
1:A:601:LEU:HD21	1:A:610:LEU:HD13	1.95	0.48
1:D:624:LEU:HB2	1:D:659:TYR:OH	2.13	0.48
1:C:617:CYS:SG	1:C:662:ARG:HD3	2.53	0.48
1:B:602:ALA:HB3	1:B:611:LEU:HD12	1.96	0.47
1:F:640:ALA:O	1:F:644:VAL:HG23	2.14	0.47
1:A:614:GLU:OE2	1:B:710:ARG:NH1	2.47	0.47
1:F:673:THR:O	1:F:675:LYS:HG3	2.14	0.47
1:C:680:ALA:O	1:C:684:MET:HG2	2.15	0.47
1:E:641:ALA:O	1:E:645:VAL:HG23	2.14	0.47
1:B:762:ARG:HB3	1:B:764:GLU:OE2	2.14	0.47
1:C:733:THR:HA	1:C:756:ASP:HB3	1.97	0.47
1:E:760:VAL:HG23	1:E:765:GLU:HB2	1.96	0.47
1:F:673:THR:O	1:F:675:LYS:NZ	2.30	0.46
1:C:657:ASP:HB3	1:C:660:GLU:OE2	2.16	0.46
1:E:736:ILE:HD13	1:E:744:LEU:HD21	1.97	0.46
1:F:723:LEU:CD2	1:F:734:VAL:HG11	2.43	0.46
1:F:744:LEU:HD23	1:F:747:ILE:HD12	1.98	0.46
1:D:655:ASN:O	1:D:658:PHE:HB3	2.16	0.46
1:D:710:ARG:HG3	1:D:712:GLN:HG3	1.97	0.46
1:F:624:LEU:HB2	1:F:659:TYR:OH	2.15	0.45
1:A:599:THR:CG2	1:A:610:LEU:HD12	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:GLY:HA3	1:A:663:ASP:CG	2.37	0.45
1:C:708:THR:OG1	1:C:710:ARG:HG2	2.16	0.45
1:C:700:ASP:OD2	1:C:700:ASP:N	2.39	0.45
1:E:680:ALA:O	1:E:684:MET:HG2	2.17	0.45
1:A:595:VAL:HA	1:A:615:THR:O	2.17	0.44
1:B:743:ASP:O	1:B:746:GLU:HG2	2.16	0.44
1:F:601:LEU:HB2	1:F:704:THR:HB	2.00	0.44
1:F:612:THR:O	1:F:669:PRO:HD2	2.18	0.44
1:B:620:GLY:HA3	1:B:663:ASP:CG	2.37	0.44
1:A:640:ALA:O	1:A:644:VAL:HG23	2.18	0.43
1:D:604:THR:CG2	1:D:611:LEU:HD21	2.48	0.43
1:B:704:THR:O	1:B:737:PRO:HD3	2.17	0.43
1:E:673:THR:HA	1:E:674:PRO:HD3	1.88	0.43
1:E:598:VAL:HG22	1:E:697:VAL:HG21	2.01	0.43
1:F:602:ALA:HB3	1:F:611:LEU:HD12	2.00	0.43
1:B:612:THR:HG21	3:C:917:HOH:O	2.19	0.43
1:D:700:ASP:OD2	1:D:700:ASP:N	2.38	0.43
1:C:723:LEU:HD12	1:C:747:ILE:CD1	2.48	0.43
1:D:748:PRO:HB2	1:D:751:VAL:HG23	2.00	0.43
1:A:599:THR:HG22	1:A:610:LEU:HD12	2.00	0.43
1:D:650:GLU:OE2	1:D:656:PRO:HG3	2.18	0.43
1:F:723:LEU:HA	1:F:723:LEU:HD23	1.69	0.42
1:E:703:MET:HE2	1:E:735:LEU:HD12	2.01	0.42
1:E:703:MET:HE3	1:E:703:MET:HB3	1.66	0.42
1:A:587:GLY:HA2	1:B:651:LYS:HG3	2.01	0.42
1:A:586:TYR:HB3	1:B:652:LEU:HD21	2.01	0.42
1:E:602:ALA:HB1	3:E:927:HOH:O	2.19	0.42
1:F:723:LEU:HD12	1:F:747:ILE:CD1	2.50	0.41
1:D:680:ALA:O	1:D:684:MET:HG2	2.20	0.41
1:D:641:ALA:O	1:D:645:VAL:HG23	2.20	0.41
1:F:626:TYR:HA	1:F:666:VAL:O	2.20	0.41
1:F:648:ARG:HH22	1:F:762:ARG:HH21	1.68	0.41
1:A:601:LEU:CD2	1:A:610:LEU:HD13	2.51	0.41
1:A:760:VAL:HG21	1:A:766:VAL:HG22	2.03	0.41
1:C:655:ASN:O	1:C:658:PHE:HB3	2.21	0.41
1:B:685:CYS:O	1:B:689:VAL:HG23	2.21	0.41
1:E:656:PRO:HD3	3:E:912:HOH:O	2.20	0.41
1:B:680:ALA:O	1:B:684:MET:HG2	2.20	0.41
1:B:735:LEU:HD12	1:B:758:HIS:HB2	2.02	0.41
1:D:659:TYR:HD2	1:D:659:TYR:HA	1.74	0.41
1:E:760:VAL:HG21	1:E:766:VAL:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:727:HIS:NE2	1:F:754:ASP:OD2	2.31	0.41
1:E:597:GLN:HE22	1:F:710:ARG:HH11	1.69	0.41
1:C:736:ILE:O	1:C:759:PRO:HA	2.20	0.40
1:E:685:CYS:O	1:E:689:VAL:HG23	2.20	0.40
1:F:736:ILE:HD13	1:F:744:LEU:HD21	2.03	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	188/200 (94%)	183 (97%)	4 (2%)	1 (0%)	29	15
1	B	179/200 (90%)	172 (96%)	7 (4%)	0	100	100
1	C	179/200 (90%)	174 (97%)	5 (3%)	0	100	100
1	D	179/200 (90%)	173 (97%)	6 (3%)	0	100	100
1	E	188/200 (94%)	182 (97%)	6 (3%)	0	100	100
1	F	177/200 (88%)	173 (98%)	4 (2%)	0	100	100
All	All	1090/1200 (91%)	1057 (97%)	32 (3%)	1 (0%)	51	37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	699	ALA

### 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	152/160 (95%)	148 (97%)	4 (3%)	46	32
1	B	145/160 (91%)	140 (97%)	5 (3%)	37	22
1	C	145/160 (91%)	141 (97%)	4 (3%)	43	29
1	D	145/160 (91%)	140 (97%)	5 (3%)	37	22
1	E	152/160 (95%)	148 (97%)	4 (3%)	46	32
1	F	143/160 (89%)	141 (99%)	2 (1%)	67	58
All	All	882/960 (92%)	858 (97%)	24 (3%)	44	30

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	591	ASN
1	A	610	LEU
1	A	659	TYR
1	A	764	GLU
1	B	612	THR
1	B	691	CYS
1	B	700	ASP
1	B	735	LEU
1	B	764	GLU
1	C	626	TYR
1	C	659	TYR
1	C	660	GLU
1	C	764	GLU
1	D	651	LYS
1	D	659	TYR
1	D	691	CYS
1	D	742	ARG
1	D	764	GLU
1	E	606	VAL
1	E	724	LEU
1	E	764	GLU
1	E	779	GLN
1	F	659	TYR
1	F	764	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	772	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	SO4	D	801	-	4,4,4	0.24	0	6,6,6	0.11	0
2	SO4	B	801	-	4,4,4	0.21	0	6,6,6	0.31	0
2	SO4	F	801	-	4,4,4	0.25	0	6,6,6	0.12	0
2	SO4	E	801	-	4,4,4	0.23	0	6,6,6	0.22	0
2	SO4	C	801	-	4,4,4	0.24	0	6,6,6	0.31	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	190/200 (95%)	-0.18	2 (1%) 80 78	24, 39, 69, 106	0
1	B	181/200 (90%)	0.07	8 (4%) 34 28	28, 45, 84, 95	0
1	C	181/200 (90%)	-0.15	2 (1%) 80 78	27, 41, 69, 90	0
1	D	181/200 (90%)	0.02	6 (3%) 46 40	33, 49, 73, 88	0
1	E	190/200 (95%)	0.18	7 (3%) 41 36	31, 51, 88, 119	0
1	F	179/200 (89%)	-0.13	1 (0%) 89 88	26, 41, 72, 98	0
All	All	1102/1200 (91%)	-0.03	26 (2%) 59 54	24, 44, 78, 119	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	659	TYR	7.4
1	E	747	ILE	5.0
1	C	606	VAL	4.5
1	A	586	TYR	4.2
1	B	751	VAL	4.0
1	E	738	PHE	3.7
1	A	775	PRO	3.4
1	E	750	ASN	3.1
1	E	752	ILE	3.0
1	E	595	VAL	2.9
1	E	753	ALA	2.9
1	D	606	VAL	2.7
1	D	775	PRO	2.6
1	B	752	ILE	2.6
1	B	606	VAL	2.6
1	D	724	LEU	2.3
1	B	742	ARG	2.3
1	D	699	ALA	2.3
1	B	608	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	751	VAL	2.2
1	F	752	ILE	2.2
1	B	775	PRO	2.2
1	D	774	GLU	2.2
1	B	700	ASP	2.1
1	C	774	GLU	2.1
1	D	728	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	SO4	F	801	5/5	0.80	0.16	98,100,101,103	0
2	SO4	C	801	5/5	0.84	0.13	57,62,66,72	0
2	SO4	E	801	5/5	0.92	0.13	62,64,66,69	0
2	SO4	B	801	5/5	0.98	0.08	43,48,51,53	0
2	SO4	D	801	5/5	0.98	0.12	54,55,58,59	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.