



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 23, 2020 – 02:04 pm GMT

PDB ID : 7AQH  
Title : Cell wall binding domain of the Staphylococcal phage 2638A endolysin  
Authors : Dunne, M.; Sobieraj, A.; Ernst, P.; Mittl, P.R.E.; Pluckthun, A.; Loessner, M.J.  
Deposited on : 2020-10-21  
Resolution : 2.49 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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.14.6  
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) : 2.14.6

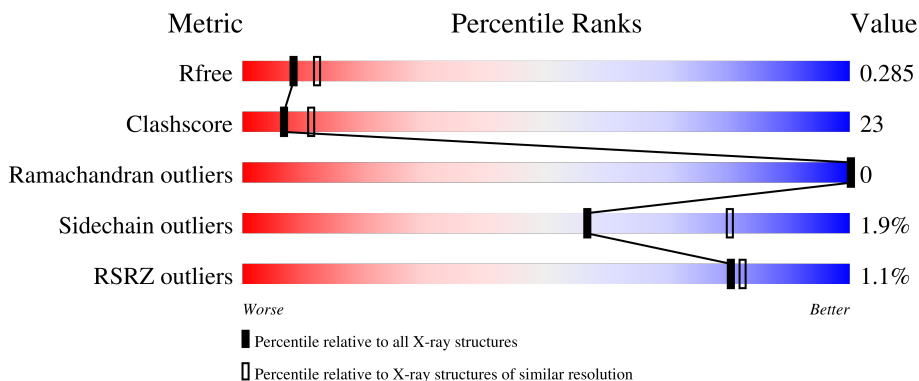
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.49 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.



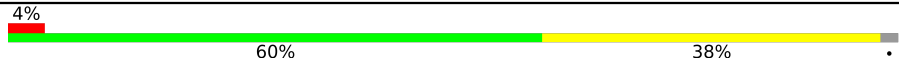


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	95	
1	B	95	
1	C	95	
1	D	95	
1	E	95	

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Mol	Chain	Length	Quality of chain
1	F	95	 4% 60% 38% .
1	G	95	 64% 33% ..
1	H	95	 55% 42% ..

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6336 atoms, of which 34 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.

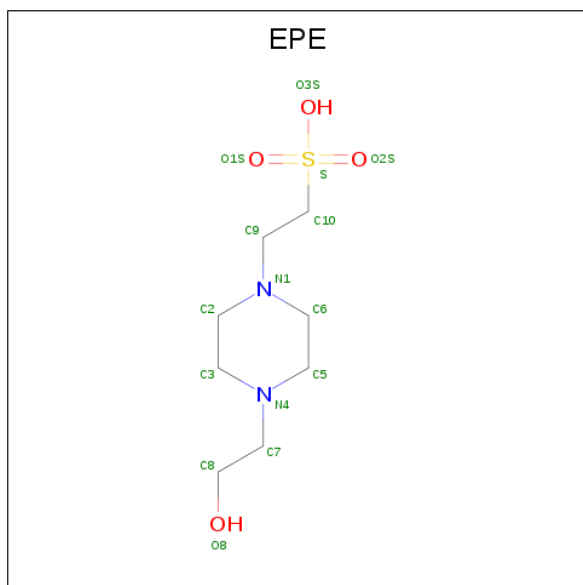
- Molecule 1 is a protein called ORF007.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	93	761	496	128	136	1	0	0	0
1	D	93	761	496	128	136	1	0	0	0
1	C	93	761	496	128	136	1	0	0	0
1	G	93	761	496	128	136	1	0	0	0
1	H	93	761	496	128	136	1	0	0	0
1	B	93	761	496	128	136	1	0	0	0
1	E	93	761	496	128	136	1	0	0	0
1	F	93	761	496	128	136	1	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	392	MET	-	initiating methionine	UNP Q4ZD58
D	392	MET	-	initiating methionine	UNP Q4ZD58
C	392	MET	-	initiating methionine	UNP Q4ZD58
G	392	MET	-	initiating methionine	UNP Q4ZD58
H	392	MET	-	initiating methionine	UNP Q4ZD58
B	392	MET	-	initiating methionine	UNP Q4ZD58
E	392	MET	-	initiating methionine	UNP Q4ZD58
F	392	MET	-	initiating methionine	UNP Q4ZD58

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C<sub>8</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf		
			Total	C	H	N	O			S	
2	A	1	Total	32	8	17	2	4	1	0	0
2	G	1	Total	32	8	17	2	4	1	0	0

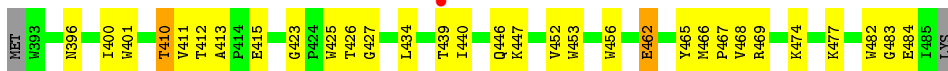
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	22	Total	22	0	0
3	D	24	Total	24	0	0
3	C	21	Total	21	0	0
3	G	22	Total	22	0	0
3	H	25	Total	25	0	0
3	B	30	Total	30	0	0
3	E	25	Total	25	0	0
3	F	15	Total	15	0	0

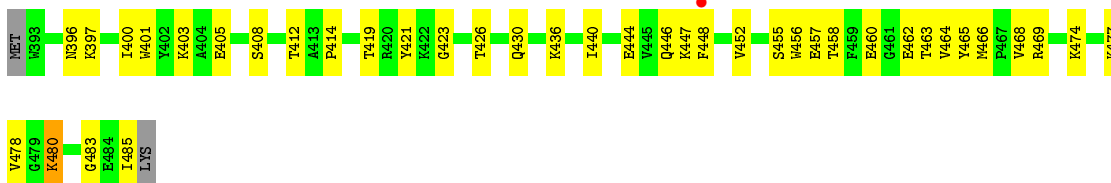
### 3 Residue-property plots [i](#)

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.

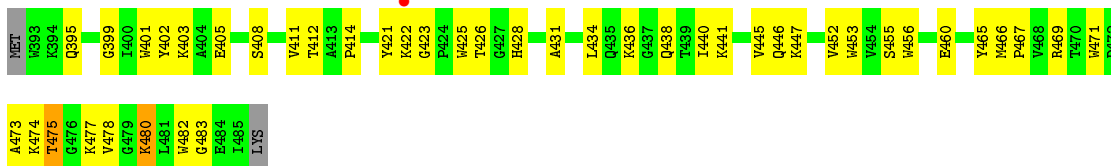
- Molecule 1: ORF007



- Molecule 1: ORF007



- Molecule 1: ORF007

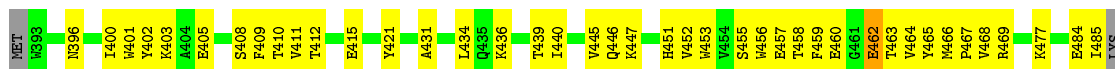


- Molecule 1: ORF007



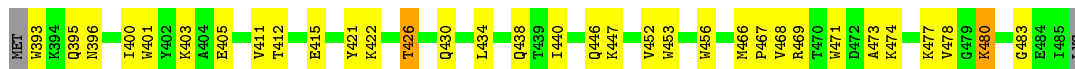
- Molecule 1: ORF007





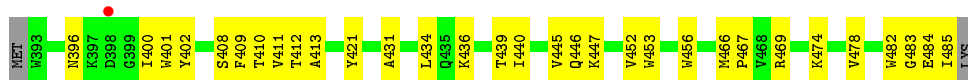
- Molecule 1: ORF007

Chain B: 63% 33%



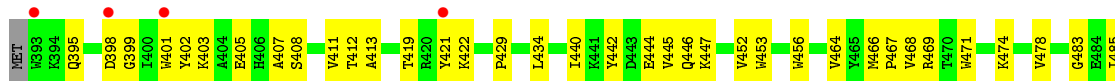
- Molecule 1: ORF007

Chain E: 65% 33%



- Molecule 1: ORF007

Chain F: 4% 60% 38%



LYS

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.13Å 62.51Å 66.17Å 111.01° 108.39° 90.18°	Depositor
Resolution (Å)	44.11 – 2.49 44.11 – 2.49	Depositor EDS
% Data completeness (in resolution range)	84.0 (44.11-2.49) 82.1 (44.11-2.49)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.06 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.15.2	Depositor
R, $R_{free}$	0.250 , 0.286 0.249 , 0.285	Depositor DCC
$R_{free}$ test set	1254 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	52.3	Xtriage
Anisotropy	0.382	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 34.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.048 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6336	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 8.56% 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: EPE

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	0.25	0/788	0.43	0/1071
1	B	0.25	0/788	0.45	0/1071
1	C	0.25	0/788	0.45	0/1071
1	D	0.27	0/788	0.45	0/1071
1	E	0.25	0/788	0.44	0/1071
1	F	0.26	0/788	0.46	0/1071
1	G	0.25	0/788	0.44	0/1071
1	H	0.24	0/788	0.44	0/1071
All	All	0.25	0/6304	0.44	0/8568

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	761	0	730	31	0
1	B	761	0	730	41	1
1	C	761	0	730	52	1
1	D	761	0	730	43	1
1	E	761	0	730	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	761	0	730	30	0
1	G	761	0	730	38	1
1	H	761	0	730	36	0
2	A	15	17	17	0	0
2	G	15	17	17	5	0
3	A	22	0	0	4	0
3	B	30	0	0	0	0
3	C	21	0	0	2	0
3	D	24	0	0	0	0
3	E	25	0	0	1	0
3	F	15	0	0	3	0
3	G	22	0	0	0	0
3	H	25	0	0	5	0
All	All	6302	34	5874	278	2

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 23.

All (278) 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:D:414:PRO:HA	1:D:436:LYS:HD2	1.26	1.07
1:D:480:LYS:HD3	1:D:480:LYS:H	1.18	1.07
1:C:480:LYS:H	1:C:480:LYS:HD2	1.21	1.02
1:B:474:LYS:HE2	1:B:474:LYS:HA	1.57	0.87
1:C:474:LYS:HA	1:C:474:LYS:HE2	1.55	0.86
1:C:414:PRO:HG2	1:B:478:VAL:HG21	1.56	0.84
1:C:480:LYS:H	1:C:480:LYS:CD	1.92	0.83
1:F:395:GLN:HE21	1:F:399:GLY:HA2	1.43	0.82
1:C:414:PRO:HG2	1:B:478:VAL:CG2	2.08	0.82
1:E:474:LYS:HA	1:E:474:LYS:HE2	1.63	0.81
1:D:458:THR:HG23	1:D:460:GLU:H	1.48	0.79
1:C:434:LEU:HD13	1:C:466:MET:HE1	1.66	0.78
1:G:396:ASN:HD21	1:G:400:ILE:HB	1.49	0.78
1:F:474:LYS:NZ	3:F:501:HOH:O	2.17	0.78
1:A:410:THR:HG23	1:A:439:THR:OG1	1.83	0.77
1:C:411:VAL:CG2	1:C:434:LEU:HD22	2.15	0.76
1:D:474:LYS:HD2	1:B:426:THR:HG23	1.67	0.76
1:B:411:VAL:HG21	1:B:434:LEU:HD22	1.68	0.76
1:C:395:GLN:HE21	1:C:399:GLY:HA2	1.51	0.76
1:D:478:VAL:HG11	1:B:415:GLU:HG3	1.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:412:THR:HG22	1:G:483:GLY:HA2	1.68	0.76
1:B:411:VAL:CG2	1:B:434:LEU:HD22	2.18	0.74
1:H:451:HIS:NE2	3:H:501:HOH:O	2.20	0.74
1:D:396:ASN:HD21	1:D:400:ILE:HB	1.53	0.73
1:A:452:VAL:HB	1:A:469:ARG:HG2	1.69	0.73
1:D:478:VAL:HG11	1:B:415:GLU:CG	2.19	0.73
1:C:411:VAL:HG21	1:C:434:LEU:HD22	1.69	0.73
1:B:480:LYS:HE2	1:B:480:LYS:H	1.54	0.72
1:G:458:THR:CG2	1:G:462:GLU:HG2	2.19	0.72
1:D:480:LYS:CD	1:D:480:LYS:H	1.98	0.72
1:G:458:THR:HG22	1:G:462:GLU:HG2	1.72	0.71
1:D:414:PRO:CA	1:D:436:LYS:HD2	2.15	0.70
1:G:480:LYS:HG3	2:G:501:EPE:H102	1.73	0.70
1:D:452:VAL:HB	1:D:469:ARG:HG2	1.72	0.70
1:H:458:THR:HG23	1:H:460:GLU:H	1.57	0.70
1:A:474:LYS:HA	1:A:474:LYS:HE2	1.72	0.70
1:D:480:LYS:N	1:D:480:LYS:HD3	2.02	0.69
1:D:457:GLU:HG2	1:D:463:THR:HG22	1.75	0.69
1:F:401:TRP:O	1:F:446:GLN:HA	1.93	0.68
1:E:452:VAL:HB	1:E:469:ARG:HG2	1.74	0.68
1:A:415:GLU:O	1:A:415:GLU:HG3	1.95	0.67
1:H:477:LYS:HD3	1:E:482:TRP:CE2	2.30	0.66
1:A:396:ASN:HD21	1:A:400:ILE:HB	1.59	0.66
1:C:434:LEU:CD2	1:C:438:GLN:HB2	2.26	0.65
1:C:401:TRP:HB2	1:C:447:LYS:HB2	1.79	0.64
1:F:452:VAL:HB	1:F:469:ARG:HG2	1.78	0.64
1:E:401:TRP:O	1:E:446:GLN:HA	1.98	0.64
1:A:412:THR:HG22	1:A:483:GLY:HA2	1.78	0.63
1:E:410:THR:HG22	1:E:439:THR:OG1	1.99	0.63
1:C:408:SER:OG	1:C:441:LYS:HG2	1.98	0.63
1:F:401:TRP:HB2	1:F:447:LYS:HB2	1.80	0.62
1:H:452:VAL:HB	1:H:469:ARG:HG2	1.81	0.62
1:B:396:ASN:HD21	1:B:400:ILE:HB	1.63	0.62
1:C:401:TRP:O	1:C:446:GLN:HA	1.98	0.62
1:G:452:VAL:HB	1:G:469:ARG:HG2	1.81	0.62
1:C:475:THR:HG22	1:C:477:LYS:H	1.63	0.62
1:H:401:TRP:O	1:H:446:GLN:HA	1.99	0.62
1:C:452:VAL:HB	1:C:469:ARG:HG2	1.81	0.61
1:H:410:THR:HG22	1:H:439:THR:OG1	2.01	0.61
1:D:396:ASN:ND2	1:D:400:ILE:HB	2.16	0.60
1:C:426:THR:CG2	1:B:474:LYS:HB3	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:426:THR:HG23	1:B:474:LYS:HB3	1.82	0.60
1:D:456:TRP:HH2	1:D:466:MET:HE3	1.66	0.60
1:C:480:LYS:N	1:C:480:LYS:HD2	2.05	0.59
1:G:396:ASN:ND2	1:G:400:ILE:HB	2.16	0.59
1:D:458:THR:HG23	1:D:460:GLU:N	2.15	0.59
1:D:421:TYR:HE1	1:D:462:GLU:HG3	1.67	0.59
1:D:456:TRP:CH2	1:D:466:MET:HE3	2.38	0.59
1:B:447:LYS:HG2	1:B:452:VAL:HG22	1.85	0.59
1:B:453:TRP:CZ3	1:B:467:PRO:HD3	2.38	0.58
1:B:434:LEU:CD2	1:B:438:GLN:HB2	2.34	0.58
1:H:458:THR:HG22	1:H:462:GLU:HG2	1.84	0.57
1:A:477:LYS:HD2	3:A:605:HOH:O	2.04	0.57
1:B:434:LEU:HD11	1:B:440:ILE:HG23	1.86	0.57
1:E:453:TRP:CZ3	1:E:467:PRO:HD3	2.39	0.57
1:B:434:LEU:HD13	1:B:466:MET:HE1	1.85	0.57
1:G:411:VAL:CG2	1:G:434:LEU:HD23	2.34	0.57
1:G:453:TRP:CZ3	1:G:467:PRO:HD3	2.40	0.57
1:G:401:TRP:O	1:G:446:GLN:HA	2.05	0.56
1:G:480:LYS:CD	1:G:480:LYS:H	2.17	0.56
1:H:453:TRP:CZ3	1:H:467:PRO:HD3	2.40	0.56
1:A:401:TRP:O	1:A:446:GLN:HA	2.05	0.56
1:C:395:GLN:HE21	1:C:399:GLY:CA	2.19	0.56
1:D:474:LYS:HD2	1:B:426:THR:CG2	2.33	0.56
1:G:480:LYS:HD2	1:G:480:LYS:H	1.71	0.56
1:A:423:GLY:HA2	1:A:465:TYR:CE1	2.41	0.56
1:C:422:LYS:HD3	1:C:428:HIS:CD2	2.40	0.56
1:C:434:LEU:HD21	1:C:438:GLN:HB2	1.88	0.55
1:C:477:LYS:HG3	3:C:514:HOH:O	2.06	0.55
1:C:412:THR:HG23	1:C:436:LYS:NZ	2.20	0.55
1:G:480:LYS:HD2	1:G:480:LYS:N	2.22	0.55
1:C:434:LEU:HD23	1:C:438:GLN:HB2	1.88	0.54
1:G:455:SER:HB3	1:G:465:TYR:CD2	2.42	0.54
1:C:434:LEU:HD11	1:C:440:ILE:HG23	1.90	0.54
1:D:458:THR:CG2	1:D:462:GLU:HG2	2.37	0.54
1:E:402:TYR:HA	1:E:445:VAL:O	2.08	0.54
1:B:480:LYS:H	1:B:480:LYS:CE	2.18	0.54
1:D:401:TRP:O	1:D:446:GLN:HA	2.07	0.54
1:C:475:THR:HG21	1:G:482:TRP:CH2	2.43	0.54
1:A:411:VAL:CG2	1:A:434:LEU:HD23	2.37	0.54
1:G:412:THR:HG22	1:G:483:GLY:CA	2.37	0.53
1:F:422:LYS:HD2	1:F:429:PRO:HD2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:477:LYS:HD2	1:G:482:TRP:CE2	2.43	0.53
1:F:452:VAL:O	1:F:468:VAL:HG22	2.08	0.53
1:E:411:VAL:CG2	1:E:434:LEU:HD23	2.39	0.53
1:D:401:TRP:HB2	1:D:447:LYS:HB2	1.89	0.52
3:A:605:HOH:O	1:F:413:ALA:HB1	2.08	0.52
1:H:477:LYS:HD3	1:E:482:TRP:CZ2	2.43	0.52
1:A:453:TRP:CZ3	1:A:467:PRO:HD3	2.45	0.52
1:A:474:LYS:NZ	3:A:602:HOH:O	2.42	0.52
1:C:395:GLN:NE2	1:C:399:GLY:HA2	2.23	0.52
1:C:412:THR:O	1:C:412:THR:HG22	2.09	0.52
1:B:480:LYS:CD	1:B:480:LYS:H	2.21	0.52
1:H:477:LYS:HE2	3:H:506:HOH:O	2.09	0.52
1:G:458:THR:HG23	1:G:460:GLU:H	1.75	0.52
1:F:402:TYR:HA	1:F:445:VAL:O	2.10	0.51
1:F:453:TRP:CZ3	1:F:467:PRO:HD3	2.45	0.51
1:D:412:THR:O	1:D:436:LYS:HE3	2.10	0.51
1:G:457:GLU:HG3	1:G:463:THR:HG22	1.91	0.51
1:D:455:SER:HB3	1:D:465:TYR:CD2	2.45	0.51
1:F:422:LYS:HE3	1:F:429:PRO:HG2	1.92	0.51
1:B:412:THR:OG1	1:B:483:GLY:HA2	2.11	0.51
1:D:400:ILE:HG12	1:D:448:PHE:CE2	2.46	0.51
1:G:458:THR:HG21	1:G:462:GLU:HG2	1.90	0.51
1:A:396:ASN:ND2	1:A:400:ILE:HB	2.25	0.51
1:G:480:LYS:HG3	2:G:501:EPE:C10	2.38	0.51
1:A:412:THR:HG22	1:A:482:TRP:O	2.11	0.51
1:H:452:VAL:O	1:H:468:VAL:HG22	2.11	0.51
3:H:505:HOH:O	1:E:413:ALA:HB1	2.11	0.50
1:C:402:TYR:HA	1:C:445:VAL:O	2.10	0.50
1:F:422:LYS:HD2	1:F:429:PRO:CD	2.42	0.50
1:H:455:SER:HB3	1:H:465:TYR:CD2	2.46	0.50
1:H:458:THR:CG2	1:H:462:GLU:HG2	2.41	0.50
1:C:434:LEU:CD1	1:C:466:MET:HE1	2.41	0.50
1:F:456:TRP:CH2	1:F:466:MET:HE3	2.47	0.50
1:H:458:THR:HG23	1:H:459:PHE:N	2.25	0.50
1:F:398:ASP:OD2	3:F:502:HOH:O	2.20	0.50
1:G:456:TRP:CH2	1:G:466:MET:HE3	2.47	0.50
1:B:396:ASN:ND2	1:B:400:ILE:HB	2.27	0.50
1:E:396:ASN:HD21	1:E:400:ILE:HB	1.76	0.50
1:D:480:LYS:HG3	2:G:501:EPE:O2S	2.11	0.49
1:G:434:LEU:HD22	1:G:466:MET:HE1	1.92	0.49
1:A:412:THR:HG22	1:A:482:TRP:C	2.32	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:403:LYS:HE3	1:H:405:GLU:OE1	2.12	0.49
1:A:456:TRP:CZ3	1:A:466:MET:HE3	2.47	0.49
1:D:412:THR:OG1	1:D:483:GLY:HA2	2.12	0.49
1:E:456:TRP:CH2	1:E:466:MET:HE3	2.46	0.49
1:C:423:GLY:HA2	1:C:465:TYR:CE1	2.47	0.49
1:H:402:TYR:HA	1:H:445:VAL:O	2.13	0.49
1:C:403:LYS:HE3	1:C:405:GLU:OE1	2.13	0.49
1:D:444:GLU:OE1	1:D:446:GLN:NE2	2.40	0.48
1:D:400:ILE:HG12	1:D:448:PHE:CZ	2.48	0.48
1:C:434:LEU:HD13	1:C:466:MET:CE	2.41	0.48
1:D:480:LYS:HD2	2:G:501:EPE:H81	1.94	0.48
1:H:447:LYS:HG2	1:H:452:VAL:HG22	1.96	0.48
1:B:474:LYS:HE2	1:B:474:LYS:CA	2.38	0.48
1:C:421:TYR:CE2	1:C:431:ALA:HA	2.49	0.48
1:E:421:TYR:CE2	1:E:431:ALA:HA	2.48	0.48
1:B:434:LEU:HD21	1:B:438:GLN:HB2	1.95	0.48
1:C:471:TRP:CZ2	1:C:473:ALA:HA	2.48	0.48
1:B:452:VAL:HB	1:B:469:ARG:HG2	1.96	0.47
1:B:456:TRP:CZ3	1:B:466:MET:HE3	2.49	0.47
1:E:411:VAL:HG23	1:E:434:LEU:HD23	1.96	0.47
1:A:462:GLU:HG2	1:A:462:GLU:H	1.40	0.47
1:F:421:TYR:O	1:F:422:LYS:HB2	2.14	0.47
1:B:434:LEU:HD23	1:B:438:GLN:HB2	1.96	0.47
1:E:434:LEU:HD22	1:E:466:MET:CE	2.45	0.47
1:A:413:ALA:HB1	3:E:505:HOH:O	2.15	0.47
1:B:471:TRP:CZ2	1:B:473:ALA:HA	2.49	0.47
1:G:408:SER:OG	1:G:441:LYS:HG2	2.15	0.47
1:C:453:TRP:CZ3	1:C:467:PRO:HD3	2.49	0.47
1:H:396:ASN:HD21	1:H:400:ILE:HB	1.80	0.47
1:F:402:TYR:HB2	1:F:444:GLU:OE2	2.14	0.47
1:H:401:TRP:HB2	1:H:447:LYS:HB2	1.97	0.47
1:F:412:THR:OG1	1:F:483:GLY:HA2	2.15	0.47
1:B:480:LYS:N	1:B:480:LYS:HE2	2.28	0.46
1:A:452:VAL:O	1:A:468:VAL:HG22	2.14	0.46
1:F:419:THR:HB	1:F:464:VAL:CG1	2.45	0.46
1:H:411:VAL:CG2	1:H:434:LEU:HD23	2.46	0.46
1:G:480:LYS:HE3	2:G:501:EPE:H101	1.98	0.46
1:H:415:GLU:HG3	1:F:478:VAL:HG11	1.97	0.46
1:D:458:THR:HG22	1:D:462:GLU:HG2	1.97	0.46
1:C:477:LYS:HD2	1:G:482:TRP:NE1	2.31	0.46
1:E:396:ASN:ND2	1:E:400:ILE:HB	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:434:LEU:HD22	1:H:466:MET:HE1	1.96	0.46
1:H:456:TRP:CZ3	1:H:466:MET:HE3	2.51	0.46
1:G:411:VAL:HG23	1:G:434:LEU:HD23	1.97	0.45
1:C:477:LYS:NZ	3:C:502:HOH:O	2.49	0.45
1:A:412:THR:HG22	1:A:483:GLY:CA	2.46	0.45
1:D:423:GLY:HA2	1:D:465:TYR:CE1	2.51	0.45
1:E:412:THR:OG1	1:E:483:GLY:HA2	2.17	0.45
1:E:434:LEU:HD22	1:E:466:MET:HE2	1.97	0.45
1:H:477:LYS:HG3	3:H:515:HOH:O	2.16	0.45
1:C:412:THR:HB	1:C:483:GLY:HA2	1.99	0.45
1:C:421:TYR:O	1:C:422:LYS:HB2	2.16	0.45
1:C:452:VAL:HG23	1:C:471:TRP:HB2	1.98	0.45
1:G:411:VAL:HG21	1:G:434:LEU:HD23	1.99	0.45
1:A:401:TRP:HB2	1:A:447:LYS:HB2	1.98	0.45
1:A:425:TRP:CH2	1:A:427:GLY:HA3	2.51	0.45
1:D:397:LYS:N	1:D:397:LYS:HD3	2.32	0.45
1:H:412:THR:HA	1:H:436:LYS:HG3	1.99	0.45
1:G:412:THR:HG22	1:G:482:TRP:C	2.38	0.45
1:C:425:TRP:HE3	1:C:428:HIS:CD2	2.35	0.44
1:A:415:GLU:CG	1:A:415:GLU:O	2.65	0.44
1:D:477:LYS:NZ	1:D:478:VAL:O	2.38	0.44
1:F:411:VAL:CG2	1:F:434:LEU:HD23	2.47	0.44
1:A:484:GLU:HA	3:A:607:HOH:O	2.17	0.44
1:E:408:SER:O	1:E:485:ILE:HA	2.18	0.44
1:F:403:LYS:HE3	1:F:405:GLU:OE1	2.18	0.44
1:D:458:THR:HG22	1:D:462:GLU:H	1.83	0.44
1:H:409:PHE:HA	1:H:484:GLU:O	2.17	0.44
1:B:401:TRP:O	1:B:446:GLN:HA	2.18	0.43
1:D:452:VAL:O	1:D:468:VAL:HG22	2.17	0.43
1:D:421:TYR:CE1	1:D:462:GLU:HG3	2.51	0.43
1:G:434:LEU:HD22	1:G:466:MET:CE	2.48	0.43
1:C:440:ILE:HD12	1:C:440:ILE:C	2.39	0.43
1:H:408:SER:O	1:H:485:ILE:HA	2.18	0.43
1:B:395:GLN:HG3	1:B:401:TRP:CZ3	2.53	0.43
1:H:440:ILE:C	1:H:440:ILE:HD12	2.39	0.43
1:B:434:LEU:CD1	1:B:466:MET:HE1	2.49	0.43
1:C:474:LYS:CA	1:C:474:LYS:HE2	2.37	0.43
1:F:440:ILE:HD12	1:F:440:ILE:C	2.38	0.43
1:A:425:TRP:CZ2	1:A:427:GLY:HA3	2.53	0.43
1:B:434:LEU:HD13	1:B:466:MET:CE	2.49	0.43
1:G:412:THR:CG2	1:G:483:GLY:HA2	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:421:TYR:O	1:B:422:LYS:HB2	2.17	0.43
1:D:478:VAL:HG11	1:B:415:GLU:HG2	1.99	0.43
1:G:401:TRP:HB2	1:G:447:LYS:HB2	1.99	0.43
1:H:457:GLU:HG2	1:H:463:THR:HG22	2.00	0.43
1:C:456:TRP:CZ3	1:C:466:MET:HE3	2.53	0.43
1:D:403:LYS:HE3	1:D:405:GLU:OE1	2.19	0.42
1:B:395:GLN:HG3	1:B:401:TRP:CE3	2.53	0.42
1:B:452:VAL:O	1:B:468:VAL:HG22	2.19	0.42
1:A:411:VAL:HG23	1:A:434:LEU:HD23	2.00	0.42
1:C:455:SER:HB3	1:C:465:TYR:CD2	2.55	0.42
1:H:455:SER:HA	1:H:464:VAL:O	2.19	0.42
1:B:403:LYS:HE3	1:B:405:GLU:OE1	2.20	0.42
1:E:440:ILE:HD12	1:E:440:ILE:C	2.40	0.42
1:E:447:LYS:HG2	1:E:452:VAL:HG22	2.00	0.42
1:C:422:LYS:HB2	1:C:428:HIS:CG	2.55	0.42
1:D:400:ILE:HG12	1:D:448:PHE:CD2	2.54	0.42
1:D:408:SER:O	1:D:485:ILE:HA	2.20	0.42
1:D:440:ILE:C	1:D:440:ILE:HD12	2.40	0.42
1:D:419:THR:HB	1:D:464:VAL:CG1	2.50	0.42
1:G:440:ILE:HD12	1:G:440:ILE:C	2.40	0.42
1:A:434:LEU:HD22	1:A:466:MET:HE1	2.02	0.42
1:C:478:VAL:HG11	1:G:415:GLU:HG3	2.01	0.42
1:B:440:ILE:HD12	1:B:440:ILE:C	2.40	0.41
1:A:440:ILE:HD12	1:A:440:ILE:C	2.40	0.41
1:B:393:TRP:CE2	1:B:403:LYS:HD3	2.56	0.41
1:G:397:LYS:N	1:G:397:LYS:HD3	2.36	0.41
1:A:447:LYS:HG2	1:A:452:VAL:HG22	2.02	0.41
1:H:415:GLU:N	3:H:503:HOH:O	2.51	0.41
1:H:421:TYR:CE2	1:H:431:ALA:HA	2.56	0.41
1:H:453:TRP:HA	1:H:466:MET:O	2.20	0.41
1:E:409:PHE:HA	1:E:484:GLU:O	2.21	0.41
1:F:447:LYS:NZ	3:F:503:HOH:O	2.40	0.41
1:H:411:VAL:HG23	1:H:434:LEU:HD23	2.03	0.41
1:F:474:LYS:CA	1:F:474:LYS:HE2	2.51	0.41
1:F:474:LYS:HA	1:F:474:LYS:HE2	2.03	0.41
1:G:402:TYR:HA	1:G:445:VAL:O	2.20	0.41
1:H:410:THR:HG22	1:H:439:THR:CG2	2.51	0.41
1:A:456:TRP:CH2	1:A:466:MET:HE3	2.56	0.41
1:C:412:THR:O	1:C:412:THR:CG2	2.67	0.41
1:C:412:THR:CB	1:C:483:GLY:HA2	2.51	0.41
1:D:397:LYS:H	1:D:397:LYS:HD3	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:398:ASP:O	1:F:398:ASP:OD1	2.39	0.41
1:A:453:TRP:HA	1:A:466:MET:O	2.20	0.40
1:F:452:VAL:HG23	1:F:471:TRP:HB2	2.04	0.40
1:G:447:LYS:HG2	1:G:452:VAL:HG22	2.02	0.40
1:G:481:LEU:HD23	1:G:481:LEU:HA	1.97	0.40
1:C:482:TRP:CE2	1:B:477:LYS:HD3	2.57	0.40
1:E:401:TRP:HB2	1:E:447:LYS:HB2	2.02	0.40
1:F:407:ALA:HB3	1:F:442:TYR:CZ	2.57	0.40
1:F:434:LEU:HD22	1:F:466:MET:CE	2.52	0.40
1:E:412:THR:HA	1:E:436:LYS:HG3	2.04	0.40
1:F:408:SER:O	1:F:485:ILE:HA	2.21	0.40
1:H:434:LEU:CD2	1:H:466:MET:HE1	2.52	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:460:GLU:OE2	1:B:430:GLN:NE2[1_655]	2.08	0.12
1:D:430:GLN:NE2	1:C:460:GLU:OE2[1_565]	2.13	0.07

## 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	91/95 (96%)	87 (96%)	4 (4%)	0	100	100
1	B	91/95 (96%)	86 (94%)	5 (6%)	0	100	100
1	C	91/95 (96%)	87 (96%)	4 (4%)	0	100	100
1	D	91/95 (96%)	87 (96%)	4 (4%)	0	100	100
1	E	91/95 (96%)	87 (96%)	4 (4%)	0	100	100
1	F	91/95 (96%)	87 (96%)	4 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	91/95 (96%)	87 (96%)	4 (4%)	0	100	100
1	H	91/95 (96%)	87 (96%)	4 (4%)	0	100	100
All	All	728/760 (96%)	695 (96%)	33 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	77/79 (98%)	74 (96%)	3 (4%)	32	57
1	B	77/79 (98%)	75 (97%)	2 (3%)	46	72
1	C	77/79 (98%)	75 (97%)	2 (3%)	46	72
1	D	77/79 (98%)	75 (97%)	2 (3%)	46	72
1	E	77/79 (98%)	76 (99%)	1 (1%)	69	87
1	F	77/79 (98%)	77 (100%)	0	100	100
1	G	77/79 (98%)	76 (99%)	1 (1%)	69	87
1	H	77/79 (98%)	76 (99%)	1 (1%)	69	87
All	All	616/632 (98%)	604 (98%)	12 (2%)	57	80

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	410	THR
1	A	426	THR
1	A	462	GLU
1	D	426	THR
1	D	480	LYS
1	C	475	THR
1	C	480	LYS
1	G	480	LYS
1	H	462	GLU

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Mol	Chain	Res	Type
1	B	426	THR
1	B	480	LYS
1	E	478	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	395	GLN
1	G	428	HIS
1	E	406	HIS
1	F	395	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](#)

2 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	EPE	G	501	-	15,15,15	0.87	1 (6%)	18,20,20	1.78	5 (27%)
2	EPE	A	501	-	15,15,15	0.84	1 (6%)	18,20,20	1.75	7 (38%)

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
2	EPE	G	501	-	-	5/9/19/19	0/1/1/1
2	EPE	A	501	-	-	3/9/19/19	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	501	EPE	C10-S	2.89	1.81	1.77
2	A	501	EPE	C10-S	2.77	1.81	1.77

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	EPE	C5-N4-C3	3.36	116.39	108.83
2	G	501	EPE	C5-N4-C3	3.33	116.33	108.83
2	G	501	EPE	C7-N4-C3	3.27	119.60	111.23
2	A	501	EPE	C7-N4-C3	3.25	119.56	111.23
2	A	501	EPE	C7-N4-C5	3.14	119.27	111.23
2	G	501	EPE	O2S-S-C10	3.03	110.56	106.92
2	G	501	EPE	C7-N4-C5	2.57	117.81	111.23
2	G	501	EPE	O1S-S-C10	2.48	109.90	106.92
2	A	501	EPE	O1S-S-C10	2.39	109.80	106.92
2	A	501	EPE	O3S-S-C10	2.17	109.28	105.77
2	A	501	EPE	C6-N1-C2	2.05	113.44	108.83
2	A	501	EPE	O2S-S-C10	2.05	109.38	106.92

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	501	EPE	S-C10-C9-N1
2	G	501	EPE	C9-C10-S-O3S
2	A	501	EPE	C10-C9-N1-C2
2	A	501	EPE	C10-C9-N1-C6
2	A	501	EPE	C8-C7-N4-C5
2	G	501	EPE	C9-C10-S-O1S
2	G	501	EPE	C9-C10-S-O2S
2	G	501	EPE	N4-C7-C8-O8

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	501	EPE	5	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	93/95 (97%)	0.18	1 (1%) 80 82	43, 51, 61, 74	0
1	B	93/95 (97%)	0.19	0 100 100	41, 50, 65, 77	0
1	C	93/95 (97%)	0.32	1 (1%) 80 82	44, 50, 60, 75	0
1	D	93/95 (97%)	0.18	1 (1%) 80 82	42, 51, 64, 76	0
1	E	93/95 (97%)	0.11	1 (1%) 80 82	41, 48, 61, 74	0
1	F	93/95 (97%)	0.33	4 (4%) 35 38	44, 53, 70, 81	0
1	G	93/95 (97%)	0.17	0 100 100	41, 51, 62, 70	0
1	H	93/95 (97%)	0.02	0 100 100	40, 50, 60, 66	0
All	All	744/760 (97%)	0.19	8 (1%) 80 82	40, 51, 63, 81	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	401	TRP	4.9
1	F	393	TRP	2.4
1	D	448	PHE	2.4
1	F	421	TYR	2.2
1	E	398	ASP	2.1
1	C	422	LYS	2.1
1	F	398	ASP	2.1
1	A	439	THR	2.1

### 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	EPE	G	501	15/15	0.41	0.28	74,93,107,117	0
2	EPE	A	501	15/15	0.75	0.20	77,92,103,103	32

### 6.5 Other polymers [i](#)

There are no such residues in this entry.