

Full wwPDB X-ray Structure Validation Report (i)

May 29, 2020 - 11:56 am BST

PDB ID	:	1WEK
Title	:	Crystal structure of the conserved hypothetical protein TT1465 from Thermus
		thermophilus HB8
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		nomics/Proteomics Initiative (RSGI)
Deposited on	:	2004-05-25
Resolution	:	2.20 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 A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.20 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	Percent	Percentile Ranks		
Clashscore			15	
W	orse	Bet	ter	
E.	Percentile relative to all X-ray stru	uctures		
01	Percentile relative to X-ray structu	ures of similar resolution		
	Whole archive	Similar reso	lution	

Metric	Whole archive	Similar resolution		
Metric	$(\# { m Entries})$	$(\# { m Entries}, { m resolution} { m range}({ m \AA}))$		
Clashscore	141614	5594(2.20-2.20)		

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain		
1	А	217	69%	27%	•
1	В	217	67%	28%	5%
1	С	217	67%	28%	5%
1	D	217	66%	30%	·
1	Е	217	67%	28%	5%
1	F	217	71%	24%	6%



$1 \mathrm{WEK}$

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 10248 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.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Δ	208	Total	С	Ν	Ο	Se	0	0	0
	Л	200	1651	1076	281	293	1	0	0	0
1	В	207	Total	С	Ν	Ο	Se	0	0	0
	D	207	1644	1071	280	292	1	0	0	0
1	С	206	Total	С	Ν	Ο	Se	0	0	0
	U	200	1632	1066	277	288	1	0	0	0
1	п	208	Total	С	Ν	Ο	Se	0	0	0
	D	200	1646	1074	279	292	1	0	0	0
1	Б	E anc	Total	С	Ν	Ο	Se	0	0	0
	200	1629	1062	276	290	1	0	0	0	
1	1 D	20.5	Total	С	Ν	Ο	Se	0	0	0
	L L	205	1624	1060	276	287	1			

• Molecule 1 is a protein called hypothetical protein TT1465.

• Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
2	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
2	С	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
2	С	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
2	С	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	Ε	1	Total O P 5 4 1	0	0
2	Ε	1	Total O P 5 4 1	0	0
2	F	1	TotalOP541	0	0
2	F	1	Total O P 5 4 1	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	75	Total O 75 75	0	0
3	В	56	Total O 56 56	0	0
3	С	48	Total O 48 48	0	0
3	D	51	$\begin{array}{cc} {\rm Total} & {\rm O} \\ 51 & 51 \end{array}$	0	0
3	Ε	65	Total O 65 65	0	0
3	F	67	Total O 67 67	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are colorcoded 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. 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.

Note EDS was not executed.



• Molecule 1: hypothetical protein TT1465



• Molecule 1: hypothetical protein TT1465





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	65.65Å 83.80 Å 265.14 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.94 - 2.20	Depositor
% Data completeness	93 9 (14 94-2 20)	Depositor
(in resolution range)	50.5 (11.51 2.20)	Depositor
R_{merge}	(Not available)	Depositor
R _{sym}	0.09	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10248	wwPDB-VP
Average B, all atoms $(Å^2)$	26.0	wwPDB-VP



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: PO4

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	
	Cham	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.42	0/1695	0.63	0/2298
1	В	0.43	0/1687	0.64	0/2287
1	С	0.41	0/1674	0.62	0/2266
1	D	0.40	0/1688	0.60	0/2286
1	Е	0.43	0/1671	0.63	0/2264
1	F	0.44	0/1666	0.61	0/2255
All	All	0.42	0/10081	0.62	0/13656

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	А	1651	0	1627	60	0
1	В	1644	0	1619	56	0
1	С	1632	0	1619	52	0
1	D	1646	0	1629	56	0
1	Е	1629	0	1605	50	0
1	F	1624	0	1608	51	0
2	А	5	0	0	0	0



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	В	5	0	0	0	0
2	С	15	0	0	1	0
2	D	15	0	0	1	0
2	Е	10	0	0	0	0
2	F	10	0	0	0	0
3	А	75	0	0	3	0
3	В	56	0	0	3	0
3	С	48	0	0	3	0
3	D	51	0	0	1	0
3	Ε	65	0	0	1	0
3	F	67	0	0	4	0
All	All	10248	0	9707	291	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 15.

All (291) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 0	Interatomic	Clash
Atom-1	Atom-2	${ m distance}~({ m \AA})$	overlap (Å)
1:C:200:THR:HG22	1:C:202:GLU:H	1.24	1.03
1:D:81:MSE:HE2	1:D:81:MSE:HA	1.52	0.88
1:F:202:GLU:HB2	1:F:205:GLU:HG3	1.55	0.87
1:A:107:LYS:HD2	1:A:107:LYS:H	1.37	0.86
1:C:159:LYS:NZ	1:D:102:LEU:HD23	1.95	0.81
1:E:30:GLU:O	1:E:34:GLU:HG2	1.81	0.81
1:F:13:GLU:HB2	3:F:519:HOH:O	1.79	0.81
1:D:157:THR:HG21	1:D:159:LYS:NZ	1.96	0.80
1:E:184:ARG:HG3	1:E:185:ASP:N	1.96	0.80
1:B:157:THR:HG23	1:B:159:LYS:HG2	1.63	0.80
1:B:61:LEU:HD22	1:B:65:LEU:HD22	1.64	0.79
1:D:151:VAL:O	1:D:155:LEU:HD23	1.82	0.79
1:B:178:ARG:HG2	1:B:178:ARG:HH11	1.48	0.78
1:A:167:LEU:HB2	1:A:199:LEU:HD13	1.63	0.78
1:E:7:ILE:HG22	1:F:130:ARG:HG2	1.67	0.77
1:C:159:LYS:HZ1	1:D:102:LEU:HD23	1.49	0.77
1:C:45:SER:HB2	1:C:140:GLY:HA2	1.66	0.76
1:A:145:LEU:CD1	1:B:149:SER:HB3	2.15	0.76
1:F:45:SER:HB2	1:F:140:GLY:HA2	1.66	0.76
1:B:204:GLU:O	1:B:208:GLN:HG2	1.86	0.75
1:F:45:SER:O	1:F:80:VAL:HG12	1.86	0.75
1:A:155:LEU:HD22	1:A:162:ARG:HA	1.69	0.74



Interstomic Clash						
Atom-1	Atom-2	distance $(Å)$	overlap(Å)			
1:F:99:ASN:HD21	1:F:112:GLN:HE22	1.36	0.73			
1:A:107:LYS:HD2	1:A:107:LYS:N	2.04	0.73			
1:D:157:THR:HG21	1:D:159:LYS:HZ3	1.53	0.72			
1:E:7:ILE:HD11	1:E:121:PHE:HE2	1.53	0.72			
1:C:202:GLU:HG3	1:C:205:GLU:HG3	1.71	0.72			
1:E:68:ALA:HB1	1:E:211:LYS:HD3	1.69	0.72			
1:D:45:SER:HB2	1:D:140:GLY:HA2	1.71	0.71			
1:A:167:LEU:HB2	1:A:199:LEU:CD1	2.20	0.70			
1:C:204:GLU:HG3	1:C:208:GLN:HE21	1.55	0.70			
1:A:45:SER:HB2	1:A:140:GLY:HA2	1.74	0.69			
1:C:101:GLU:C	1:C:102:LEU:HD12	2.13	0.68			
1:E:200:THR:HG22	1:E:201:ASP:N	2.07	0.68			
1:C:122:PHE:HA	1:C:125:LYS:HD3	1.76	0.68			
1:B:44:GLY:HA3	1:B:80:VAL:CG2	2.25	0.67			
1:D:99:ASN:ND2	1:D:108:PRO:HB3	2.10	0.67			
1:E:42:VAL:HG11	1:E:84:VAL:HG21	1.77	0.67			
1:A:113:THR:HG22	1:A:114:HIS:CD2	2.30	0.66			
1:D:42:VAL:HG11	1:D:84:VAL:HG21	1.77	0.66			
1:C:149:SER:HB3	1:D:146:ASP:HA	1.78	0.66			
2:C:507:PO4:O3	1:D:12:HIS:HD2	1.79	0.66			
1:B:162:ARG:HD2	3:B:512:HOH:O	1.94	0.66			
1:E:45:SER:HB2	1:E:140:GLY:HA2	1.78	0.65			
1:E:167:LEU:HB2	1:E:199:LEU:HD22	1.79	0.65			
1:C:204:GLU:O	1:C:208:GLN:HG2	1.95	0.65			
1:C:138:LEU:HD21	1:C:168:LEU:HD12	1.78	0.64			
1:C:26:VAL:O	1:C:30:GLU:HG3	1.96	0.64			
1:E:99:ASN:ND2	1:E:108:PRO:HB3	2.13	0.64			
1:A:161:HIS:HB3	1:A:163:PHE:CE1	2.31	0.64			
1:A:98:LEU:O	1:A:124:ARG:HD2	1.97	0.64			
1:E:65:LEU:HD13	1:E:207:VAL:HG22	1.79	0.64			
1:A:155:LEU:CD2	1:A:162:ARG:HA	2.28	0.63			
1:B:161:HIS:ND1	1:B:162:ARG:O	2.30	0.63			
1:B:157:THR:CG2	1:B:159:LYS:HG2	2.27	0.63			
1:D:81:MSE:CE	1:D:81:MSE:HA	2.27	0.63			
1:E:146:ASP:HA	1:F:149:SER:HB3	1.81	0.62			
1:D:157:THR:HG22	1:D:159:LYS:HG3	1.82	0.62			
1:D:20:ARG:O	1:D:24:GLU:HG3	2.01	0.61			
1:A:12:HIS:ND1	1:B:33:SER:HB2	2.15	0.61			
1:E:200:THR:HG22	1:E:202:GLU:H	1.65	0.61			
1:F:7:ILE:HD11	1:F:121:PHE:HZ	1.66	0.61			
1:F:162:ARG:HG3	1:F:162:ARG:HH11	1.64	0.61			



A 4 1	A 4 5 775 D	Interatomic	Clash
Atom-1	Atom-2	${ m distance}~({ m \AA})$	overlap (Å)
1:A:107:LYS:CD	1:A:107:LYS:H	2.13	0.60
1:A:145:LEU:CD1	1:B:149:SER:CB	2.80	0.60
1:E:189:VAL:HG13	1:E:193:ASP:HB2	1.84	0.60
1:D:162:ARG:HD3	1:D:162:ARG:O	2.02	0.60
1:A:61:LEU:HD11	1:A:136:VAL:HG11	1.82	0.59
1:A:177:VAL:HG21	1:A:199:LEU:HD22	1.84	0.59
1:E:200:THR:HG22	1:E:201:ASP:H	1.65	0.59
1:C:82:GLU:HG3	1:C:109:ASN:HB2	1.85	0.59
1:B:202:GLU:HB3	1:B:205:GLU:HG3	1.85	0.59
1:D:125:LYS:HD2	1:D:150:GLU:HG3	1.84	0.59
1:E:7:ILE:CG2	1:F:130:ARG:HG2	2.33	0.59
1:A:125:LYS:HD3	1:A:147:GLU:OE1	2.02	0.59
1:C:189:VAL:HG13	1:C:193:ASP:HB2	1.85	0.59
1:E:125:LYS:HD3	1:E:150:GLU:OE1	2.04	0.58
1:F:74:THR:HB	1:F:81:MSE:HG3	1.84	0.58
1:A:202:GLU:HB3	1:A:204:GLU:CD	2.23	0.58
1:A:22:LEU:HD11	1:B:22:LEU:HD21	1.84	0.58
1:C:137:PHE:HE2	1:C:147:GLU:HB3	1.69	0.57
1:F:81:MSE:SE	3:F:553:HOH:O	2.71	0.57
1:C:18:LEU:HD23	1:D:22:LEU:HD22	1.86	0.57
1:D:189:VAL:HG13	1:D:193:ASP:HB2	1.87	0.57
1:F:125:LYS:HD3	1:F:150:GLU:OE1	2.05	0.57
1:F:7:ILE:HD11	1:F:121:PHE:CZ	2.39	0.57
1:E:160:VAL:HG13	1:F:7:ILE:HD12	1.87	0.56
1:B:63:ARG:O	1:B:67:GLU:HG3	2.04	0.56
1:C:181:ALA:O	1:C:184:ARG:HG2	2.04	0.56
1:E:86:ARG:O	1:E:90:GLU:HG3	2.05	0.56
1:B:44:GLY:HA3	1:B:80:VAL:HG21	1.86	0.56
1:C:80:VAL:O	1:C:84:VAL:HG23	2.06	0.56
1:F:207:VAL:HG12	1:F:211:LYS:HD2	1.88	0.56
1:B:44:GLY:HA3	1:B:80:VAL:HG22	1.88	0.56
1:D:173:TRP:O	1:D:177:VAL:HG13	2.05	0.56
1:E:179:TRP:O	1:E:182:PHE:HB3	2.05	0.56
1:F:100:ILE:HG12	1:F:124:ARG:NH1	2.21	0.56
1:D:137:PHE:HB2	1:D:167:LEU:HD23	1.87	0.56
1:A:6:LEU:H	1:A:6:LEU:HD23	1.71	0.55
1:D:155:LEU:HD21	1:D:163:PHE:CZ	2.41	0.55
1:E:117:SER:C	1:E:118:LEU:HD12	2.27	0.55
1:C:138:LEU:HD23	1:C:168:LEU:HB2	1.89	0.55
1:B:161:HIS:ND1	1:B:162:ARG:N	2.55	0.55
1:C:176:LEU:O	1:C:180:LEU:HG	2.07	0.55



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:178:ARG:NH1	1:B:178:ARG:HG2	2.17	0.55
1:B:80:VAL:O	1:B:84:VAL:HG23	2.07	0.55
1:D:213:GLU:HG2	1:D:214:ALA:N	2.21	0.54
1:E:7:ILE:HD11	1:E:121:PHE:CE2	2.38	0.54
1:A:26:VAL:O	1:A:30:GLU:HG3	2.06	0.54
1:D:192:GLU:O	1:D:195:GLN:HB2	2.06	0.54
1:C:182:PHE:CE1	1:C:186:GLN:NE2	2.75	0.54
1:D:157:THR:HG21	1:D:159:LYS:HZ2	1.72	0.54
1:D:179:TRP:O	1:D:182:PHE:HB3	2.07	0.54
1:B:202:GLU:HG3	1:B:204:GLU:OE1	2.09	0.53
1:B:211:LYS:O	1:B:211:LYS:HG2	2.08	0.53
1:F:177:VAL:HG21	1:F:199:LEU:HD11	1.90	0.53
1:A:121:PHE:HE1	1:B:159:LYS:HG3	1.73	0.53
1:B:161:HIS:HD1	1:B:162:ARG:N	2.07	0.53
1:A:80:VAL:O	1:A:84:VAL:HG23	2.09	0.52
1:B:167:LEU:HB2	1:B:199:LEU:HD13	1.90	0.52
1:E:109:ASN:HD22	1:E:111:TYR:H	1.57	0.52
1:C:55:TYR:CD1	1:C:83:ALA:HB2	2.44	0.52
1:A:121:PHE:CE1	1:B:159:LYS:HG3	2.45	0.52
1:C:116:LEU:N	1:C:116:LEU:HD23	2.24	0.52
1:D:153:VAL:O	1:D:157:THR:HB	2.09	0.52
1:D:100:ILE:HD11	1:D:124:ARG:HD2	1.90	0.52
1:D:186:GLN:O	1:D:187:LYS:HB2	2.09	0.52
1:A:101:GLU:OE2	1:A:119:ARG:HG2	2.09	0.52
1:A:95:SER:H	1:A:113:THR:HB	1.75	0.52
1:F:38:PRO:HB2	1:F:70:PHE:HD2	1.74	0.52
1:A:184:ARG:HA	1:A:189:VAL:HG12	1.91	0.52
1:B:162:ARG:HG3	1:B:163:PHE:N	2.25	0.52
1:D:176:LEU:HD13	1:D:176:LEU:C	2.30	0.52
1:F:9:GLN:HE22	1:F:119:ARG:HH22	1.57	0.51
1:B:168:LEU:O	1:B:169:ASP:HB3	2.11	0.51
1:C:177:VAL:HG21	1:C:199:LEU:HD21	1.92	0.51
1:B:45:SER:HB2	1:B:140:GLY:HA2	1.92	0.51
1:E:200:THR:CG2	1:E:201:ASP:N	2.74	0.51
1:A:207:VAL:O	1:A:211:LYS:HG3	2.11	0.51
1:B:78:PRO:HG2	1:B:106:GLN:HG2	1.93	0.51
1:C:164:PRO:HB2	1:C:166:PHE:CE1	2.46	0.51
1:F:179:TRP:O	1:F:182:PHE:HB3	2.11	0.50
1:B:204:GLU:CD	1:B:204:GLU:H	2.14	0.50
1:C:184:ARG:HB2	3:C:550:HOH:O	2.10	0.50
1:F:162:ARG:HG3	1:F:162:ARG:NH1	2.26	0.50



		Interatomic	Clash	
Atom-1	Atom-2	distance $(Å)$	overlap (Å)	
1:A:20:ARG:NH2	1:F:27:GLU:HG2	2.26	0.50	
1:B:86:ARG:HD2	1:B:111:TYR:CE1	2.46	0.50	
1:C:147:GLU:O	1:C:151:VAL:HG23	2.11	0.50	
1:F:38:PRO:HB2	1:F:70:PHE:CD2	2.46	0.50	
1:E:4:LYS:HE3	1:E:8:ASP:HB3	1.93	0.50	
1:A:161:HIS:HD2	1:A:163:PHE:CE2	2.30	0.50	
1:F:119:ARG:NH1	3:F:511:HOH:O	2.44	0.50	
1:C:149:SER:CB	1:D:146:ASP:HA	2.42	0.49	
1:C:179:TRP:O	1:C:182:PHE:HB3	2.12	0.49	
1:D:157:THR:CG2	1:D:157:THR:O	2.60	0.49	
1:F:86:ARG:O	1:F:90:GLU:HG3	2.12	0.49	
1:D:16:TRP:CZ3	1:F:13:GLU:HG3	2.47	0.49	
1:E:153:VAL:HG21	1:F:146:ASP:HB2	1.94	0.49	
1:F:64:ALA:HB1	1:F:207:VAL:HG21	1.94	0.49	
1:B:97:GLY:HA3	1:B:112:GLN:NE2	2.28	0.49	
1:D:101:GLU:OE1	1:D:119:ARG:NE	2.45	0.49	
1:B:210:LEU:C	1:B:212:ALA:H	2.15	0.49	
1:D:6:LEU:HD11	1:D:119:ARG:O	2.13	0.49	
1:E:214:ALA:H	1:E:215:PRO:CD	2.26	0.49	
1:B:52:HIS:ND1	1:B:53:PRO:HD2	2.28	0.49	
1:A:80:VAL:CG1	1:A:138:LEU:HD13	2.43	0.48	
1:A:179:TRP:CD1	1:B:175:GLY:HA3	2.48	0.48	
1:E:200:THR:CG2	1:E:201:ASP:H	2.27	0.48	
1:C:142:PHE:CE2	1:D:189:VAL:HG21	2.49	0.48	
1:C:15:SER:O	1:C:18:LEU:HB2	2.12	0.48	
1:C:146:ASP:HA	1:D:149:SER:HB3	1.96	0.48	
1:A:189:VAL:HG22	1:A:193:ASP:HB2	1.96	0.47	
1:E:168:LEU:O	1:E:169:ASP:CB	2.62	0.47	
1:A:189:VAL:HG22	1:A:193:ASP:OD2	2.14	0.47	
1:E:133:VAL:O	1:E:164:PRO:HD2	2.13	0.47	
2:D:512:PO4:O4	1:E:30:GLU:OE1	2.32	0.47	
1:A:105:GLU:HA	1:A:105:GLU:OE1	2.14	0.47	
1:C:200:THR:HG23	1:C:205:GLU:OE1	2.13	0.47	
1:C:43:PHE:CD1	1:C:147:GLU:HG3	2.50	0.47	
1:A:204:GLU:HG2	1:A:205:GLU:N	2.30	0.47	
1:A:179:TRP:CZ3	1:B:176:LEU:HG	2.50	0.47	
1:A:95:SER:O	1:A:113:THR:N	2.47	0.46	
1:F:125:LYS:HD3	1:F:150:GLU:CD	2.35	0.46	
1:A:9:GLN:O	1:A:13:GLU:HG3	2.16	0.46	
1:F:186:GLN:O	1:F:187:LYS:HB2	2.15	0.46	
1:C:186:GLN:O	1:C:187:LYS:HB2	2.15	0.46	



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:F:26:VAL:O	1:F:30:GLU:HG3	2.16	0.46
1:F:44:GLY:HA3	1:F:80:VAL:HG11	1.97	0.46
1:B:20:ARG:O	1:B:24:GLU:HG3	2.16	0.46
1:C:18:LEU:HD23	1:D:22:LEU:CD2	2.46	0.46
1:F:169:ASP:HA	1:F:201:ASP:OD1	2.16	0.46
1:B:60:ARG:HD2	3:B:547:HOH:O	2.16	0.46
1:F:80:VAL:O	1:F:84:VAL:HG23	2.15	0.46
1:F:9:GLN:NE2	1:F:119:ARG:HH22	2.14	0.45
1:E:200:THR:HG22	1:E:202:GLU:N	2.32	0.45
1:B:18:LEU:O	1:B:22:LEU:HB2	2.16	0.45
1:A:145:LEU:HD12	1:B:149:SER:CB	2.46	0.45
1:A:22:LEU:HD11	1:B:22:LEU:CD2	2.46	0.45
1:B:65:LEU:HD13	1:B:207:VAL:HG22	1.97	0.45
1:E:125:LYS:HD3	1:E:150:GLU:CD	2.36	0.45
1:F:202:GLU:CB	1:F:205:GLU:HG3	2.38	0.45
1:C:200:THR:HG22	1:C:201:ASP:N	2.32	0.44
1:A:168:LEU:O	1:A:169:ASP:HB3	2.16	0.44
1:C:31:THR:HG21	1:C:116:LEU:CD2	2.47	0.44
1:C:174:GLU:HG2	1:C:178:ARG:NH2	2.33	0.44
1:D:157:THR:O	1:D:158:GLU:HB2	2.18	0.44
1:A:129:VAL:HG23	3:A:563:HOH:O	2.17	0.44
1:D:80:VAL:CG1	1:D:138:LEU:HD13	2.47	0.44
1:B:86:ARG:HB2	1:B:111:TYR:CD1	2.53	0.44
1:B:119:ARG:HG2	1:B:120:TYR:CE1	2.52	0.44
1:C:102:LEU:HD12	1:C:102:LEU:N	2.31	0.44
1:C:174:GLU:HG2	1:C:178:ARG:HH22	1.83	0.44
1:E:65:LEU:HD23	1:E:72:VAL:HG12	2.00	0.44
1:A:20:ARG:HB3	3:A:520:HOH:O	2.18	0.44
1:B:125:LYS:HG2	1:B:147:GLU:OE2	2.18	0.44
1:C:168:LEU:O	1:C:169:ASP:CB	2.66	0.44
1:E:137:PHE:CZ	1:E:148:LEU:HA	2.53	0.44
1:F:204:GLU:O	1:F:208:GLN:HG3	2.17	0.44
1:E:153:VAL:HG21	1:F:146:ASP:CB	2.48	0.43
1:D:5:PRO:HA	3:D:519:HOH:O	2.17	0.43
1:E:176:LEU:O	1:E:180:LEU:HG	2.18	0.43
1:F:162:ARG:NH2	3:F:529:HOH:O	2.51	0.43
1:E:175:GLY:HA3	1:F:179:TRP:CD1	2.53	0.43
1:A:101:GLU:HG3	1:A:118:LEU:O	2.18	0.43
1:B:55:TYR:CE1	1:B:83:ALA:HB2	2.53	0.43
1:E:160:VAL:HG13	1:F:7:ILE:CD1	2.48	0.43
1:C:162:ARG:NH1	1:C:162:ARG:HG3	2.34	0.43



Interatomic Clash							
Atom-1	Atom-2	distance $(Å)$	overlan (Å)				
1:E:177:VAL:HG21	1:E:199:LEU:HD23	2.00	0.43				
1:B:82:GLU:HG3	1:B:109:ASN:HB2	2.01	0.43				
1:D:168:LEU:HA	1:D:200:THR:O	2.19	0.43				
1:E:186:GLN:O	1:E:187:LYS:HB2	2.19	0.43				
1:A:20:ARG:NH1	1:A:119:ARG:HG3	2.34	0.43				
1:A:55:TYR:CE1	1:A:83:ALA:HB2	2.53	0.43				
1:A:20:ARG:HD2	3:A:520:HOH:O	2.17	0.43				
1:D:12:HIS:CE1	1:D:13:GLU:HG3	2.54	0.43				
1:D:121:PHE:O	1:D:125:LYS:HG2	2.19	0.42				
3:E:554:HOH:O	1:F:125:LYS:HD2	2.19	0.42				
1:F:17:ARG:NH1	1:F:20:ARG:HH21	2.17	0.42				
1:C:202:GLU:HA	1:C:203:PRO:HD3	1.87	0.42				
1:F:164:PRO:HG2	1:F:213:GLU:OE1	2.19	0.42				
1:B:211:LYS:HE2	1:B:211:LYS:HB3	1.86	0.42				
1:F:168:LEU:O	1:F:169:ASP:CB	2.67	0.42				
1:A:145:LEU:HD12	1:B:149:SER:HB3	1.97	0.42				
1:C:8:ASP:OD2	1:D:161:HIS:ND1	2.48	0.42				
1:D:82:GLU:HG3	1:D:109:ASN:HB2	2.01	0.42				
1:E:109:ASN:HA	1:E:110:PRO:HD3	1.84	0.42				
1:F:208:GLN:HE21	1:F:208:GLN:HB3	1.63	0.42				
1:F:209:ALA:O	1:F:213:GLU:HG3	2.19	0.42				
1:B:107:LYS:HA	1:B:108:PRO:HD3	1.85	0.42				
1:B:63:ARG:NH1	3:B:545:HOH:O	2.52	0.42				
1:C:160:VAL:HG22	1:C:161:HIS:N	2.35	0.42				
1:C:162:ARG:HG3	1:C:162:ARG:HH11	1.83	0.42				
1:A:174:GLU:HG3	1:A:178:ARG:HE	1.85	0.42				
1:A:190:GLY:N	1:A:193:ASP:OD2	2.50	0.42				
1:C:168:LEU:O	1:C:169:ASP:HB3	2.19	0.42				
3:C:513:HOH:O	1:D:150:GLU:HG2	2.19	0.42				
1:A:174:GLU:O	1:A:178:ARG:HG3	2.19	0.42				
1:A:19:PHE:CE2	1:E:18:LEU:HD12	2.54	0.42				
1:E:6:LEU:HD11	1:E:119:ARG:O	2.20	0.42				
1:F:112:GLN:NE2	1:F:115:ALA:HA	2.34	0.42				
1:B:194:LEU:HD12	1:B:194:LEU:HA	1.85	0.42				
1:B:102:LEU:O	1:B:105:GLU:HG2	2.20	0.41				
1:A:30:GLU:O	1:A:34:GLU:HG3	2.19	0.41				
1:B:116:LEU:N	1:B:116:LEU:HD12	2.35	0.41				
1:A:189:VAL:HG22	1:A:193:ASP:CB	2.51	0.41				
1:D:168:LEU:HD11	1:D:206:VAL:HG21	2.03	0.41				
1:E:4:LYS:HA	1:E:5:PRO:HD3	1.87	0.41				
1:C:168:LEU:HA	1:C:200:THR:O	2.20	0.41				



A 4 1	A 4 5 55 0	Interatomic	Clash
Atom-1	Atom-2	${ m distance}~({ m \AA})$	overlap (Å)
1:C:99:ASN:ND2	1:C:108:PRO:HB3	2.35	0.41
1:C:5:PRO:HB3	1:D:158:GLU:O	2.20	0.41
1:D:170:ARG:HG2	1:D:199:LEU:HB3	2.02	0.41
1:D:34:GLU:OE2	1:E:119:ARG:NH2	2.45	0.41
1:D:214:ALA:HA	1:D:215:PRO:HD3	1.82	0.41
1:D:64:ALA:HB1	1:D:207:VAL:HG21	2.02	0.41
1:A:80:VAL:HG11	1:A:138:LEU:HD13	2.03	0.41
1:B:66:ALA:HB1	1:B:93:GLY:HA3	2.03	0.41
1:D:20:ARG:HD3	1:E:30:GLU:OE2	2.20	0.41
1:E:100:ILE:HB	1:E:124:ARG:NH2	2.36	0.41
1:F:125:LYS:HD3	1:F:150:GLU:OE2	2.20	0.41
1:D:148:LEU:HD12	1:D:173:TRP:CZ3	2.56	0.40
1:A:107:LYS:CD	1:A:107:LYS:N	2.74	0.40
1:C:32:LEU:HB3	1:C:131:TYR:CD1	2.57	0.40
1:A:55:TYR:CD1	1:A:83:ALA:HB2	2.57	0.40
1:B:202:GLU:HA	1:B:203:PRO:HD3	2.01	0.40
1:D:181:ALA:HA	1:D:194:LEU:HD21	2.03	0.40
1:E:168:LEU:O	1:E:169:ASP:HB3	2.22	0.40
1:E:65:LEU:CD1	1:E:207:VAL:HG22	2.50	0.40
1:A:138:LEU:O	1:A:139:PRO:C	2.57	0.40
1:A:16:TRP:CH2	1:A:17:ARG:CZ	3.04	0.40
3:C:513:HOH:O	1:D:125:LYS:NZ	2.55	0.40
1:F:155:LEU:HD22	1:F:162:ARG:HD3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains (i)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

12 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).

Mal	Tuno	Chain	Dec	Tink	B	ond leng	\mathbf{gths}	E	ond ang	gles
	туре	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
2	PO4	C	507	-	4,4,4	1.62	0	6,6,6	0.45	0
2	PO4	В	503	-	4,4,4	1.52	0	$6, \! 6, \! 6$	0.44	0
2	PO4	F	502	-	4,4,4	1.62	0	$6,\!6,\!6$	0.44	0
2	PO4	D	506	-	4,4,4	1.55	0	$6, \! 6, \! 6$	0.46	0
2	PO4	Е	509	-	4,4,4	1.55	0	$6,\!6,\!6$	0.41	0
2	PO4	Е	504	-	4,4,4	1.39	0	$6,\!6,\!6$	0.43	0
2	PO4	А	501	-	4,4,4	1.61	0	$6, \! 6, \! 6$	0.43	0
2	PO4	С	511	-	4,4,4	1.70	0	6,6,6	0.44	0
2	PO4	D	508	-	4,4,4	1.66	1 (25%)	$6,\!6,\!6$	0.46	0
2	PO4	D	512	-	4,4,4	1.55	0	6,6,6	0.39	0
2	PO4	С	505	-	4,4,4	1.70	0	6,6,6	0.47	0
2	PO4	F	510	-	4,4,4	1.58	0	$6, \! 6, \! 6$	0.43	0

All (1) bond length outliers are listed below:

Mol	Chain	\mathbf{Res}	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(\operatorname{\AA})$
2	D	508	PO4	P-04	-2.03	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.



There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	С	507	PO4	1	0
2	D	512	PO4	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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

