



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 3, 2023 – 10:18 pm GMT

PDB ID : 2WJZ  
Title : Crystal structure of (HisH) K181A Y138A mutant of imidazoleglycerolphosphate synthase (HisH HisF) which displays constitutive glutaminase activity  
Authors : Vega, M.C.; List, F.; Razeto, A.; Haeger, M.C.; Babinger, K.; Kuper, J.; Sterner, R.; Wilmanns, M.  
Deposited on : 2009-06-02  
Resolution : 2.60 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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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

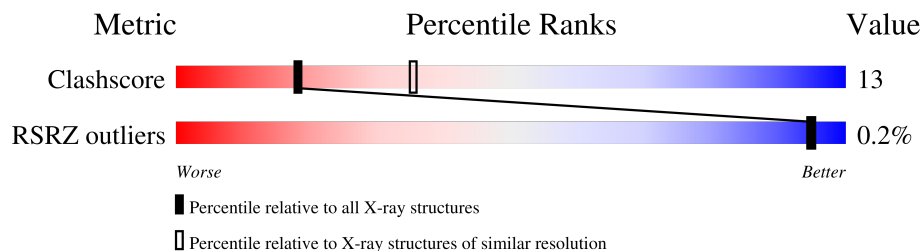
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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))
Clashscore	141614	3518 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	253	
1	C	253	
1	E	253	
2	B	201	
2	D	201	
2	F	201	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10203 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 IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE HISF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	251	1883	1200	317	360	6	0	0	0
1	C	251	1888	1200	320	362	6	0	0	0
1	E	237	1769	1128	298	337	6	0	0	0

- Molecule 2 is a protein called IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE SUB-UNIT HISH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	193	1474	942	250	274	8	0	0	0
2	D	196	1525	978	258	281	8	0	0	0
2	F	196	1440	920	245	267	8	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	181	ALA	LYS	engineered mutation	UNP Q9X0C8
B	138	ALA	TYR	engineered mutation	UNP Q9X0C8
D	181	ALA	LYS	engineered mutation	UNP Q9X0C8
D	138	ALA	TYR	engineered mutation	UNP Q9X0C8
F	181	ALA	LYS	engineered mutation	UNP Q9X0C8
F	138	ALA	TYR	engineered mutation	UNP Q9X0C8

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0
3	E	1	Total O P 5 4 1	0	0

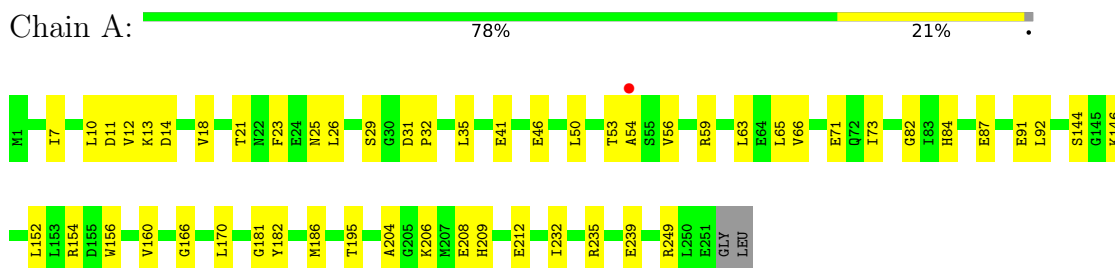
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	43	Total O 43 43	0	0
4	B	31	Total O 31 31	0	0
4	C	59	Total O 59 59	0	0
4	D	36	Total O 36 36	0	0
4	E	32	Total O 32 32	0	0
4	F	8	Total O 8 8	0	0

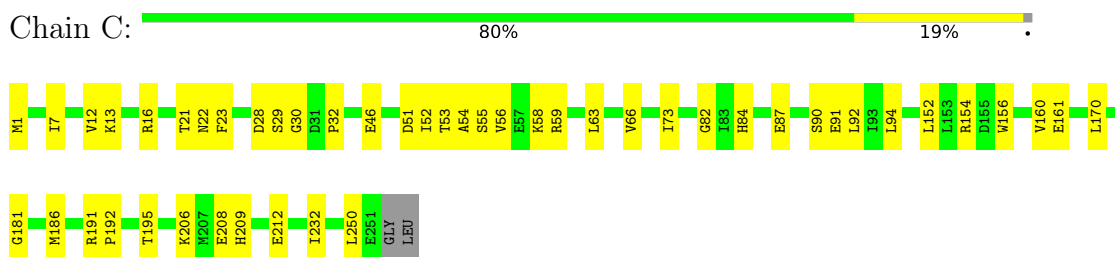
### 3 Residue-property plots

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.

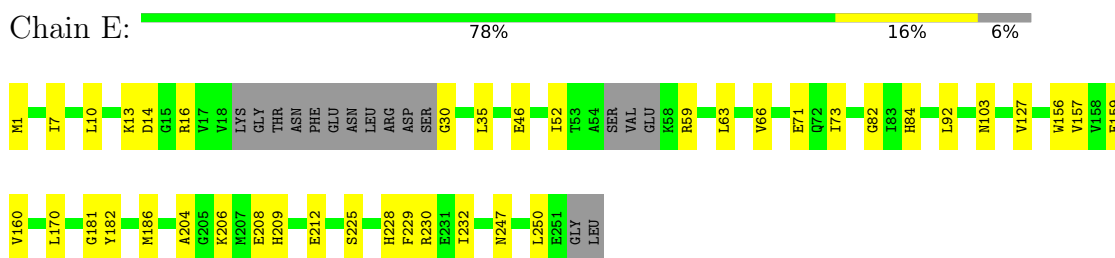
- Molecule 1: IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE HISF



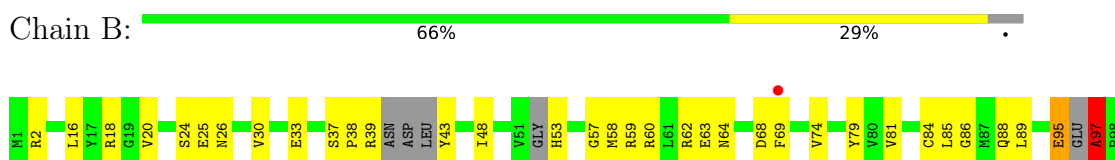
- Molecule 1: IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE HISF

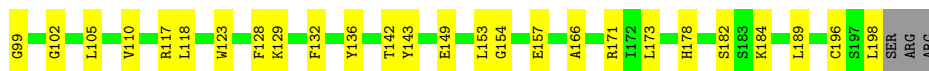


- Molecule 1: IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE HISF



- Molecule 2: IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE SUBUNIT HISH





- Molecule 2: IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE SUBUNIT HISH

Chain D: 67% 31%



- Molecule 2: IMIDAZOLE GLYCEROL PHOSPHATE SYNTHASE SUBUNIT HISH

Chain F: 70% 27%



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.94Å 93.94Å 166.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.51 – 2.60 39.51 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.51-2.60) 99.7 (39.51-2.60)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.01 (at 2.61Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.185 , 0.218 0.185 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	52.8	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 54.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l 0.119 for h,-h-k,-l 0.039 for -k,-h,-l	Xtriage
Reported twinning fraction	0.125 for h,-h-k,-l	Depositor
Outliers	0 of 50261 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10203	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% 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: 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	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.31	0/1910	0.49	0/2582
1	C	0.30	0/1915	0.49	0/2588
1	E	0.31	0/1792	0.49	0/2423
2	B	0.38	1/1504 (0.1%)	0.59	3/2026 (0.1%)
2	D	0.33	0/1557	0.52	1/2102 (0.0%)
2	F	0.31	0/1472	0.61	2/1988 (0.1%)
All	All	0.32	1/10150 (0.0%)	0.53	6/13709 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	95	GLU	CA-C	-7.85	1.32	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	8	VAL	CB-CA-C	-14.50	83.85	111.40
2	B	99	GLY	N-CA-C	8.07	133.27	113.10
2	B	97	ALA	CB-CA-C	7.18	120.87	110.10
2	B	95	GLU	CB-CA-C	-6.46	97.47	110.40
2	F	8	VAL	N-CA-C	6.41	128.31	111.00
2	D	99	GLY	N-CA-C	5.31	126.37	113.10



There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	97	ALA	Mainchain

## 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	1883	0	1872	44	0
1	C	1888	0	1880	52	0
1	E	1769	0	1756	34	0
2	B	1474	0	1398	50	0
2	D	1525	0	1477	51	0
2	F	1440	0	1335	45	0
3	A	5	0	0	0	0
3	C	5	0	0	1	0
3	E	5	0	0	1	0
4	A	43	0	0	2	0
4	B	31	0	0	6	0
4	C	59	0	0	3	0
4	D	36	0	0	9	0
4	E	32	0	0	8	0
4	F	8	0	0	2	0
All	All	10203	0	9718	265	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 13.

All (265) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:197:SER:O	2:F:198:LEU:HD23	1.34	1.26
2:B:43:TYR:N	4:B:2009:HOH:O	1.80	1.12
1:C:59:ARG:HG3	1:C:59:ARG:HH11	1.15	1.10
1:A:59:ARG:HG3	1:A:59:ARG:HH11	1.13	1.05
1:E:59:ARG:HG3	1:E:59:ARG:HH11	1.14	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:42:LEU:N	4:D:2009:HOH:O	1.93	1.02
1:E:160:VAL:N	4:E:2024:HOH:O	1.96	0.97
2:D:149:GLU:OE2	4:D:2024:HOH:O	1.83	0.96
1:C:21:THR:HG22	1:C:23:PHE:H	1.34	0.89
1:E:206:LYS:HA	1:E:232:ILE:HD11	1.52	0.89
2:F:8:VAL:HG23	2:F:9:GLY:N	1.89	0.86
1:C:59:ARG:HG3	1:C:59:ARG:NH1	1.90	0.85
1:C:21:THR:HB	1:C:29:SER:O	1.78	0.83
1:E:157:VAL:C	4:E:2024:HOH:O	2.17	0.82
1:A:59:ARG:HG3	1:A:59:ARG:NH1	1.88	0.81
1:E:59:ARG:HG3	1:E:59:ARG:NH1	1.89	0.80
1:C:16:ARG:HG2	1:C:16:ARG:HH11	1.47	0.80
2:F:95:GLU:C	2:F:97:ALA:N	2.33	0.77
2:F:42:LEU:HD12	2:F:78:ARG:HH21	1.49	0.77
1:C:16:ARG:HH11	1:C:16:ARG:CG	1.98	0.76
1:A:21:THR:CG2	1:A:23:PHE:HD1	1.98	0.75
1:A:21:THR:HG22	1:A:23:PHE:H	1.51	0.75
2:B:68:ASP:OD2	1:C:13:LYS:NZ	2.19	0.74
1:E:156:TRP:O	4:E:2024:HOH:O	2.05	0.74
2:B:95:GLU:C	2:B:97:ALA:N	2.41	0.74
2:B:53:HIS:HB3	4:B:2012:HOH:O	1.87	0.73
1:C:13:LYS:HG3	1:C:52:ILE:HG13	1.71	0.73
2:F:95:GLU:C	2:F:97:ALA:H	1.91	0.73
2:B:63:GLU:HG2	1:C:16:ARG:CZ	2.19	0.72
1:C:152:LEU:HD13	1:C:154:ARG:NH2	2.04	0.72
1:C:21:THR:CG2	1:C:23:PHE:HD1	2.03	0.72
1:C:16:ARG:HG2	1:C:16:ARG:NH1	2.03	0.71
1:A:206:LYS:HD2	1:A:208:GLU:OE1	1.90	0.71
2:B:38:PRO:HA	2:B:69:PHE:CE1	2.25	0.71
1:C:206:LYS:HD2	1:C:208:GLU:OE1	1.90	0.71
2:B:63:GLU:HA	1:C:16:ARG:HD2	1.73	0.71
1:A:10:LEU:HD13	1:A:35:LEU:HD23	1.73	0.71
2:D:107:GLU:O	4:D:2016:HOH:O	2.08	0.70
1:A:152:LEU:HD13	1:A:154:ARG:NH2	2.06	0.70
3:C:1252:PO4:O4	4:C:2059:HOH:O	2.10	0.69
1:C:206:LYS:HA	1:C:232:ILE:HD11	1.75	0.68
2:D:95:GLU:C	2:D:97:ALA:N	2.41	0.68
1:A:166:GLY:O	4:A:2025:HOH:O	2.11	0.67
2:B:26:ASN:ND2	4:B:2007:HOH:O	2.28	0.67
2:F:8:VAL:CG2	2:F:9:GLY:N	2.56	0.66
2:D:92:GLU:OE1	4:D:2015:HOH:O	2.14	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:157:VAL:O	4:E:2024:HOH:O	2.13	0.65
1:E:159:GLU:N	4:E:2024:HOH:O	2.29	0.65
1:C:21:THR:HG23	1:C:23:PHE:CD1	2.32	0.64
1:A:14:ASP:OD1	1:A:53:THR:HG23	1.97	0.64
2:D:20:VAL:HG12	2:D:32:ILE:HD13	1.80	0.64
2:F:56:GLU:HG2	2:F:59:ARG:HH12	1.63	0.64
2:F:197:SER:C	2:F:198:LEU:HD23	2.16	0.63
1:A:249:ARG:HG3	2:B:136:TYR:HB2	1.81	0.63
2:D:95:GLU:C	2:D:97:ALA:H	2.00	0.63
1:C:21:THR:CG2	1:C:23:PHE:CD1	2.82	0.63
2:B:38:PRO:C	2:B:39:ARG:HD2	2.20	0.62
1:E:206:LYS:HD2	1:E:208:GLU:OE1	1.98	0.62
1:C:195:THR:HB	2:D:117:ARG:HH11	1.63	0.62
1:A:21:THR:HG23	1:A:23:PHE:HD1	1.64	0.61
1:C:87:GLU:O	1:C:91:GLU:HG2	2.01	0.61
1:A:21:THR:HG23	1:A:23:PHE:CD1	2.36	0.60
2:D:58:MET:SD	2:D:102:GLY:HA2	2.41	0.60
1:A:87:GLU:O	1:A:91:GLU:HG2	2.02	0.60
2:F:7:SER:HB3	2:F:34:LEU:HD22	1.83	0.60
1:E:1:MET:N	4:E:2001:HOH:O	2.35	0.60
2:D:6:ILE:O	4:D:2001:HOH:O	2.17	0.60
2:B:25:GLU:HG2	4:B:2006:HOH:O	2.02	0.60
1:C:208:GLU:O	1:C:212:GLU:HG3	2.03	0.59
1:A:21:THR:CG2	1:A:23:PHE:CD1	2.82	0.59
2:D:53:HIS:CD2	2:D:96:GLU:HG2	2.38	0.59
1:C:21:THR:HG21	1:C:23:PHE:HD1	1.66	0.58
2:D:26:ASN:ND2	4:D:2007:HOH:O	2.36	0.58
1:A:59:ARG:NH2	4:A:2007:HOH:O	1.94	0.58
1:C:53:THR:O	1:C:58:LYS:HB2	2.03	0.58
1:E:16:ARG:NH2	1:E:30:GLY:HA3	2.19	0.58
2:B:58:MET:SD	2:B:102:GLY:HA2	2.44	0.57
1:C:195:THR:HB	2:D:117:ARG:NH1	2.18	0.57
1:E:7:ILE:HG12	1:E:46:GLU:HB2	1.86	0.57
2:B:24:SER:HB2	2:B:30:VAL:HG21	1.85	0.56
1:C:7:ILE:HG12	1:C:46:GLU:HB2	1.87	0.56
2:F:58:MET:SD	2:F:102:GLY:HA2	2.45	0.56
1:C:181:GLY:HA2	1:C:209:HIS:CG	2.40	0.56
2:F:6:ILE:HA	2:F:35:VAL:HB	1.87	0.56
2:B:74:VAL:HG21	2:B:105:LEU:HD22	1.88	0.55
1:C:21:THR:CG2	1:C:22:ASN:N	2.69	0.55
1:E:181:GLY:HA2	1:E:209:HIS:CG	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:LEU:CD1	1:A:35:LEU:HD23	2.36	0.55
1:A:13:LYS:O	1:A:14:ASP:HB2	2.07	0.55
1:A:41:GLU:O	2:B:184:LYS:HE2	2.07	0.55
1:E:10:LEU:HD13	1:E:35:LEU:HD23	1.87	0.55
2:B:2:ARG:NH1	2:B:33:GLU:OE2	2.39	0.55
1:E:71:GLU:HA	2:F:18:ARG:HD3	1.88	0.55
1:A:181:GLY:HA2	1:A:209:HIS:CG	2.42	0.55
1:A:208:GLU:O	1:A:212:GLU:HG3	2.07	0.55
2:F:1:MET:HE2	2:F:198:LEU:HD21	1.88	0.55
1:A:7:ILE:HG12	1:A:46:GLU:HB2	1.88	0.54
2:F:74:VAL:HG21	2:F:105:LEU:HD22	1.89	0.54
2:F:1:MET:CE	2:F:198:LEU:HD21	2.38	0.54
2:D:74:VAL:HG21	2:D:105:LEU:HD22	1.88	0.54
2:B:2:ARG:NH2	2:B:39:ARG:HB2	2.23	0.54
2:F:53:HIS:CD2	2:F:96:GLU:HG2	2.43	0.54
1:A:195:THR:HB	2:B:117:ARG:HH11	1.73	0.53
1:E:103:ASN:HB3	3:E:1252:PO4:O1	2.09	0.53
2:F:128:PHE:HD2	2:F:153:LEU:HD22	1.73	0.53
2:B:2:ARG:HH22	2:B:39:ARG:HB2	1.72	0.53
2:B:85:LEU:O	2:B:89:LEU:HD13	2.09	0.53
1:C:21:THR:HG23	1:C:23:PHE:HD1	1.71	0.53
1:C:55:SER:HA	1:C:58:LYS:HB3	1.89	0.53
2:F:85:LEU:O	2:F:89:LEU:HD13	2.09	0.52
2:D:39:ARG:O	2:D:39:ARG:HG3	2.09	0.52
2:D:74:VAL:HG21	2:D:105:LEU:CD2	2.39	0.52
1:C:12:VAL:HA	1:C:16:ARG:O	2.10	0.52
2:B:59:ARG:O	2:B:63:GLU:HG3	2.09	0.52
2:D:35:VAL:HG21	2:D:43:TYR:CE1	2.45	0.52
2:F:95:GLU:O	2:F:97:ALA:N	2.41	0.52
1:C:30:GLY:O	1:C:32:PRO:HD3	2.10	0.52
1:C:21:THR:CG2	1:C:23:PHE:H	2.16	0.52
1:A:23:PHE:HB2	1:A:26:LEU:HD12	1.92	0.52
1:E:208:GLU:O	1:E:212:GLU:HG3	2.10	0.52
2:B:97:ALA:N	4:B:2018:HOH:O	2.42	0.51
2:D:128:PHE:HD2	2:D:153:LEU:HD22	1.75	0.51
2:F:74:VAL:HG21	2:F:105:LEU:CD2	2.41	0.51
1:A:206:LYS:HA	1:A:232:ILE:HD11	1.93	0.51
1:E:13:LYS:HB2	1:E:52:ILE:HG13	1.92	0.51
1:A:73:ILE:HD13	1:A:73:ILE:O	2.11	0.51
2:B:95:GLU:O	2:B:97:ALA:N	2.44	0.51
2:F:42:LEU:HB2	4:F:2003:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:59:ARG:NH1	1:E:59:ARG:CG	2.66	0.51
2:F:197:SER:O	2:F:198:LEU:CD2	2.30	0.50
2:D:43:TYR:HB2	2:D:73:HIS:NE2	2.26	0.50
2:D:59:ARG:O	2:D:63:GLU:HG3	2.11	0.50
2:B:128:PHE:HD2	2:B:153:LEU:HD22	1.76	0.50
2:B:25:GLU:CG	4:B:2006:HOH:O	2.58	0.50
2:D:95:GLU:O	2:D:97:ALA:N	2.44	0.50
1:C:63:LEU:HD23	1:C:92:LEU:HD23	1.94	0.50
1:C:16:ARG:NH2	1:C:28:ASP:OD1	2.45	0.50
2:F:81:VAL:HG22	2:F:173:LEU:HB2	1.94	0.50
1:A:195:THR:HB	2:B:117:ARG:NH1	2.26	0.49
1:A:63:LEU:HD23	1:A:92:LEU:HD23	1.95	0.49
2:B:74:VAL:HG21	2:B:105:LEU:CD2	2.42	0.49
1:A:59:ARG:NH1	1:A:59:ARG:CG	2.66	0.49
2:D:157:GLU:HG3	2:D:157:GLU:O	2.11	0.49
2:D:118:LEU:HD13	2:D:142:THR:HG23	1.94	0.49
1:C:55:SER:HA	1:C:58:LYS:CB	2.43	0.48
1:A:144:SER:HB2	1:A:146:LYS:HE3	1.96	0.48
1:E:13:LYS:O	1:E:14:ASP:HB2	2.13	0.48
1:A:12:VAL:HG21	1:A:65:LEU:HD22	1.95	0.48
1:A:21:THR:HB	1:A:29:SER:O	2.14	0.48
2:B:37:SER:OG	2:B:38:PRO:HD2	2.14	0.48
2:D:79:TYR:CE1	2:D:171:ARG:HD3	2.48	0.47
2:D:85:LEU:O	2:D:89:LEU:HD13	2.14	0.47
1:E:63:LEU:HD23	1:E:92:LEU:HD23	1.96	0.47
2:F:157:GLU:O	2:F:157:GLU:HG3	2.14	0.47
1:C:73:ILE:HD13	1:C:73:ILE:O	2.14	0.47
2:D:26:ASN:ND2	4:D:2006:HOH:O	2.46	0.47
2:D:130:ASP:OD2	2:D:168:ARG:NH2	2.46	0.47
1:E:73:ILE:HD13	1:E:73:ILE:O	2.14	0.47
1:A:66:VAL:HG21	1:A:92:LEU:HD22	1.97	0.47
2:F:129:LYS:NZ	2:F:149:GLU:O	2.47	0.47
2:D:48:ILE:O	2:D:86:GLY:HA3	2.15	0.47
1:C:13:LYS:CG	1:C:52:ILE:HG13	2.44	0.47
2:D:57:GLY:O	2:D:61:LEU:HG	2.14	0.47
1:C:16:ARG:HH22	1:C:28:ASP:CG	2.19	0.46
1:E:66:VAL:HG21	1:E:92:LEU:HD22	1.96	0.46
2:B:117:ARG:O	2:B:118:LEU:HD23	2.15	0.46
2:D:129:LYS:NZ	2:D:149:GLU:O	2.49	0.46
1:A:21:THR:HG21	1:A:23:PHE:HD1	1.75	0.46
2:B:38:PRO:HA	2:B:69:PHE:CD1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:79:TYR:CE1	2:B:171:ARG:HD3	2.51	0.46
1:C:195:THR:CB	2:D:117:ARG:NH1	2.79	0.46
2:F:95:GLU:O	2:F:96:GLU:C	2.52	0.46
2:D:97:ALA:O	2:D:98:PRO:C	2.53	0.46
1:A:11:ASP:O	1:A:18:VAL:HG23	2.16	0.46
2:B:129:LYS:NZ	2:B:149:GLU:O	2.49	0.46
1:C:161:GLU:OE1	4:C:2036:HOH:O	2.21	0.45
2:B:48:ILE:O	2:B:86:GLY:HA3	2.17	0.45
2:D:95:GLU:O	2:D:96:GLU:C	2.53	0.45
1:C:66:VAL:HG21	1:C:92:LEU:HD22	1.97	0.45
2:F:118:LEU:HD13	2:F:142:THR:HG23	1.98	0.45
2:F:130:ASP:OD2	2:F:168:ARG:NH2	2.49	0.45
2:B:16:LEU:HD22	2:B:20:VAL:HG23	1.97	0.45
2:B:118:LEU:HD13	2:B:142:THR:HG23	1.98	0.45
2:B:81:VAL:HG22	2:B:173:LEU:HB2	1.98	0.45
2:D:81:VAL:HG22	2:D:173:LEU:HB2	1.98	0.45
2:F:16:LEU:HD22	2:F:20:VAL:HG23	1.98	0.45
2:B:57:GLY:H	2:B:60:ARG:HG3	1.81	0.45
2:D:39:ARG:NH1	2:D:39:ARG:HG2	2.31	0.45
2:F:58:MET:HA	2:F:58:MET:CE	2.47	0.44
2:F:56:GLU:CG	2:F:59:ARG:HH12	2.28	0.44
2:F:57:GLY:H	2:F:60:ARG:HG3	1.82	0.44
2:D:57:GLY:H	2:D:60:ARG:HG3	1.83	0.44
2:B:157:GLU:HG3	2:B:157:GLU:O	2.15	0.44
1:C:12:VAL:HG23	1:C:12:VAL:O	2.17	0.44
1:C:156:TRP:O	1:C:160:VAL:HG23	2.18	0.44
2:D:1:MET:HE2	2:D:198:LEU:HD21	1.99	0.44
1:C:12:VAL:O	1:C:51:ASP:HA	2.18	0.44
1:C:170:LEU:HD21	1:C:186:MET:HG2	1.98	0.44
2:B:88:GLN:HB3	2:B:110:VAL:HG21	2.00	0.44
2:F:37:SER:HB2	2:F:38:PRO:HD2	1.98	0.44
1:C:1:MET:N	4:C:2001:HOH:O	2.52	0.43
1:E:1:MET:HG3	1:E:247:ASN:ND2	2.34	0.43
1:C:90:SER:O	1:C:94:LEU:HG	2.18	0.43
1:A:170:LEU:HD21	1:A:186:MET:HG2	2.01	0.43
1:A:182:TYR:CZ	1:A:204:ALA:HB2	2.53	0.43
2:D:151:HIS:HD2	4:D:2016:HOH:O	2.01	0.43
2:F:13:ILE:C	2:F:13:ILE:HD12	2.39	0.43
2:F:42:LEU:N	4:F:2003:HOH:O	2.51	0.43
1:C:54:ALA:O	1:C:56:VAL:N	2.52	0.43
2:D:13:ILE:C	2:D:13:ILE:HD12	2.40	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:58:MET:HA	2:D:58:MET:CE	2.49	0.43
2:D:84:CYS:SG	2:D:178:HIS:CE1	3.12	0.43
2:D:123:TRP:CH2	2:D:182:SER:HA	2.54	0.42
2:B:84:CYS:SG	2:B:178:HIS:CE1	3.12	0.42
2:B:110:VAL:HG12	2:B:143:TYR:HB3	2.01	0.42
1:C:250:LEU:HA	1:C:250:LEU:HD23	1.82	0.42
2:D:38:PRO:O	2:D:39:ARG:HB3	2.19	0.42
2:D:16:LEU:HD22	2:D:20:VAL:HG23	2.00	0.42
1:E:228:HIS:C	1:E:230:ARG:H	2.22	0.42
2:F:48:ILE:O	2:F:86:GLY:HA3	2.20	0.42
1:A:23:PHE:HB3	1:A:25:ASN:OD1	2.20	0.42
1:A:54:ALA:O	1:A:56:VAL:N	2.53	0.42
1:A:71:GLU:HA	2:B:18:ARG:HD3	2.02	0.42
2:B:58:MET:HA	2:B:58:MET:CE	2.50	0.42
1:E:182:TYR:CZ	1:E:204:ALA:HB2	2.55	0.42
2:B:123:TRP:CH2	2:B:182:SER:HA	2.54	0.42
1:A:31:ASP:HA	1:A:32:PRO:HD2	1.84	0.42
1:A:235:ARG:O	1:A:239:GLU:HG2	2.20	0.42
1:E:82:GLY:O	1:E:84:HIS:HD2	2.02	0.42
2:F:117:ARG:O	2:F:118:LEU:HD23	2.20	0.42
1:C:82:GLY:O	1:C:84:HIS:HD2	2.02	0.41
2:D:88:GLN:HB3	2:D:110:VAL:HG21	2.02	0.41
1:E:230:ARG:NH2	4:E:2029:HOH:O	2.54	0.41
2:F:84:CYS:SG	2:F:178:HIS:CE1	3.13	0.41
2:B:110:VAL:CG1	2:B:143:TYR:HB3	2.50	0.41
1:C:59:ARG:NH1	1:C:59:ARG:CG	2.66	0.41
2:F:62:ARG:C	2:F:64:ASN:H	2.24	0.41
2:F:110:VAL:CG1	2:F:143:TYR:HB3	2.50	0.41
2:B:196:CYS:C	2:B:198:LEU:H	2.24	0.41
2:D:155:THR:CG2	4:D:2012:HOH:O	2.68	0.41
2:D:62:ARG:C	2:D:64:ASN:H	2.24	0.41
1:E:16:ARG:HG2	4:E:2007:HOH:O	2.21	0.41
2:B:154:GLY:N	2:B:166:ALA:HB3	2.36	0.41
1:C:21:THR:HG22	1:C:22:ASN:N	2.35	0.41
2:D:196:CYS:C	2:D:198:LEU:H	2.23	0.41
1:A:156:TRP:O	1:A:160:VAL:HG23	2.21	0.41
2:B:37:SER:O	2:B:39:ARG:HD2	2.21	0.41
1:E:225:SER:O	1:E:229:PHE:N	2.48	0.41
1:A:82:GLY:O	1:A:84:HIS:HD2	2.04	0.41
2:D:15:ASN:HB3	2:D:180:GLU:HG3	2.03	0.41
2:F:96:GLU:HB2	2:F:143:TYR:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:110:VAL:HG12	2:F:143:TYR:HB3	2.02	0.41
2:F:132:PHE:CD2	2:F:189:LEU:HD22	2.56	0.41
1:C:191:ARG:HB3	1:C:192:PRO:HD3	2.02	0.41
2:F:32:ILE:H	2:F:32:ILE:HD12	1.86	0.41
2:B:62:ARG:C	2:B:64:ASN:H	2.24	0.40
2:D:84:CYS:HG	2:D:178:HIS:CE1	2.39	0.40
1:E:156:TRP:O	1:E:160:VAL:HG23	2.20	0.40
2:B:132:PHE:CD2	2:B:189:LEU:HD22	2.56	0.40
2:D:39:ARG:O	2:D:39:ARG:CG	2.65	0.40
1:E:170:LEU:HD21	1:E:186:MET:HG2	2.03	0.40
1:E:127:VAL:HG21	1:E:160:VAL:HG13	2.03	0.40
2:F:123:TRP:CH2	2:F:182:SER:HA	2.56	0.40
1:A:50:LEU:HD12	1:A:50:LEU:HA	1.94	0.40
1:E:250:LEU:HD23	1:E:250:LEU:HA	1.78	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 monosaccharides in this entry.



## 5.6 Ligand geometry

3 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
3	PO4	A	1252	-	4,4,4	2.36	3 (75%)	6,6,6	0.52	0
3	PO4	E	1252	-	4,4,4	2.40	3 (75%)	6,6,6	0.56	0
3	PO4	C	1252	-	4,4,4	2.43	3 (75%)	6,6,6	0.43	0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1252	PO4	P-O3	-2.68	1.46	1.54
3	C	1252	PO4	P-O3	-2.67	1.46	1.54
3	E	1252	PO4	P-O3	-2.66	1.46	1.54
3	C	1252	PO4	P-O2	-2.66	1.46	1.54
3	C	1252	PO4	P-O4	-2.62	1.46	1.54
3	E	1252	PO4	P-O2	-2.60	1.46	1.54
3	E	1252	PO4	P-O4	-2.55	1.46	1.54
3	A	1252	PO4	P-O2	-2.53	1.47	1.54
3	A	1252	PO4	P-O4	-2.48	1.47	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
3	E	1252	PO4	1	0
3	C	1252	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](#)

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	251/253 (99%)	-0.31	1 (0%) 92   91	45, 68, 103, 125	0
1	C	251/253 (99%)	-0.31	0 100   100	43, 66, 100, 117	0
1	E	237/253 (93%)	-0.33	0 100   100	47, 69, 103, 137	0
2	B	193/201 (96%)	-0.17	1 (0%) 91   89	49, 76, 111, 140	0
2	D	196/201 (97%)	-0.12	1 (0%) 91   89	47, 74, 112, 137	0
2	F	196/201 (97%)	-0.19	0 100   100	53, 78, 118, 145	0
All	All	1324/1362 (97%)	-0.25	3 (0%) 95   95	43, 71, 108, 145	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	69	PHE	4.0
2	D	1	MET	2.7
1	A	54	ALA	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( $\text{\AA}^2$ )	Q<0.9
3	PO4	E	1252	5/5	0.95	0.20	80,94,124,138	0
3	PO4	A	1252	5/5	0.97	0.16	81,98,121,128	0
3	PO4	C	1252	5/5	0.98	0.15	81,86,102,107	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.