



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:57 AM GMT

PDB ID : 2YW3
Title : Crystal Structure Analysis of the 4-hydroxy-2-oxoglutarate aldolase/2-dehydro-3-deoxyphosphogluconate aldolase from TTHB1
Authors : Kawano, Y.; Hashimoto, K.; Kamiya, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-04-19
Resolution : 1.67 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

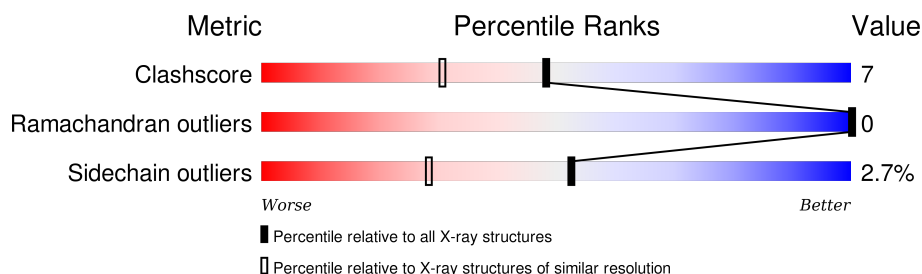
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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	102246	5317 (1.70-1.66)
Ramachandran outliers	100387	5225 (1.70-1.66)
Sidechain outliers	100360	5224 (1.70-1.66)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	207	
1	B	207	
1	C	207	
1	D	207	
1	E	207	
1	F	207	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	B	2006	-	-	X	-
2	PO4	D	2012	-	-	X	-
2	PO4	E	2014	-	-	X	-

2 Entry composition [i](#)

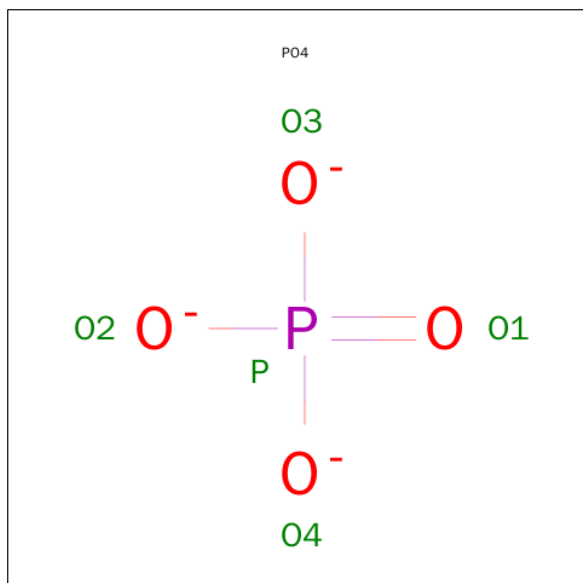
There are 3 unique types of molecules in this entry. The entry contains 10084 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 4-hydroxy-2-oxoglutarate aldolase/2-deydro-3-deoxyphosphogluconate aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O		0	3	0
			1465	948	256	261				
1	B	197	Total	C	N	O		0	3	0
			1480	958	258	264				
1	C	197	Total	C	N	O		0	4	0
			1488	964	257	267				
1	D	199	Total	C	N	O	S	0	10	0
			1529	996	259	273	1			
1	E	200	Total	C	N	O	S	0	5	0
			1509	980	260	268	1			
1	F	199	Total	C	N	O	S	0	5	0
			1506	978	259	268	1			

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



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	132	Total O 132 132	0	0
3	B	132	Total O 132 132	0	0
3	C	130	Total O 130 130	0	0

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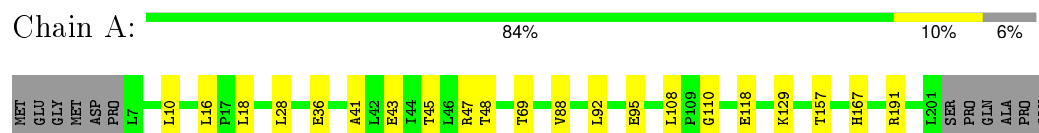
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	220	Total 220	O 220	0	0
3	E	198	Total 198	O 198	0	0
3	F	211	Total 211	O 211	0	0

3 Residue-property plots [i](#)

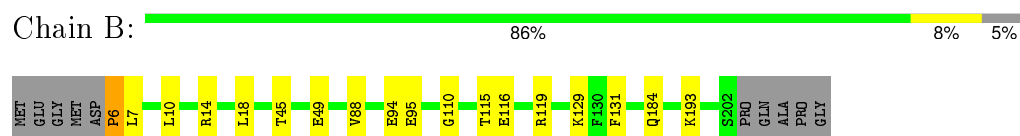
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

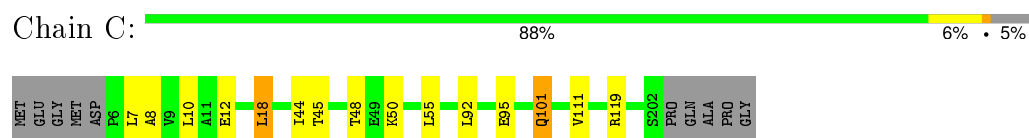
- Molecule 1: 4-hydroxy-2-oxoglutarate aldolase/2-deydro-3-deoxyphosphogluconate aldolase



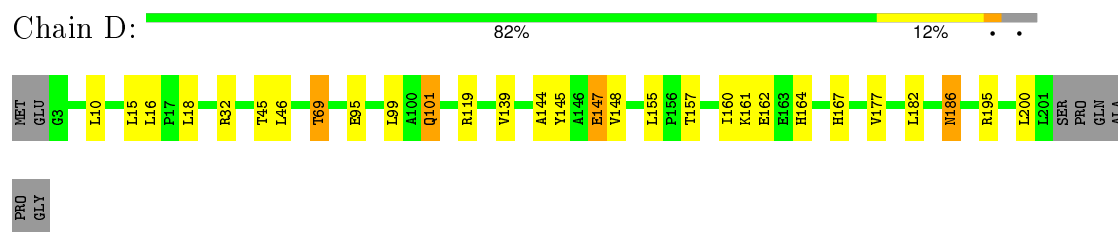
- Molecule 1: 4-hydroxy-2-oxoglutarate aldolase/2-deydro-3-deoxyphosphogluconate aldolase



- Molecule 1: 4-hydroxy-2-oxoglutarate aldolase/2-deydro-3-deoxyphosphogluconate aldolase



- Molecule 1: 4-hydroxy-2-oxoglutarate aldolase/2-deydro-3-deoxyphosphogluconate aldolase



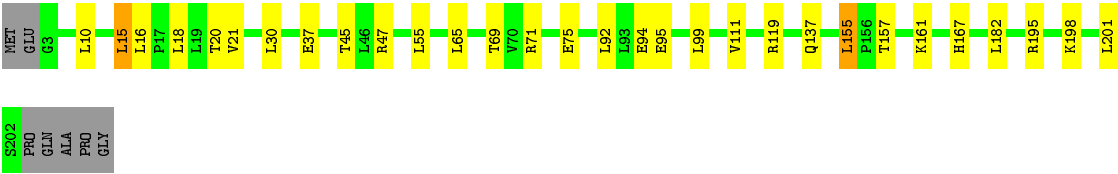
- Molecule 1: 4-hydroxy-2-oxoglutarate aldolase/2-deydro-3-deoxyphosphogluconate aldolase

Chain E:

82%

14%

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• Molecule 1: 4-hydroxy-2-oxoglutarate aldolase/2-deydro-3-deoxyphosphogluconate aldolase

Chain F:

83%

12%

••



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	125.92Å 72.69Å 161.31Å 90.00° 91.07° 90.00°	Depositor
Resolution (Å)	35.58 – 1.67	Depositor
% Data completeness (in resolution range)	100.0 (35.58-1.67)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.182 , 0.214	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10084	wwPDB-VP
Average B, all atoms (Å ²)	19.0	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	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.60	0/1498	0.65	0/2033
1	B	0.60	0/1513	0.65	1/2053 (0.0%)
1	C	0.62	0/1525	0.64	0/2070
1	D	0.74	0/1584	0.80	1/2153 (0.0%)
1	E	0.73	0/1549	0.78	2/2104 (0.1%)
1	F	0.72	0/1546	0.75	0/2100
All	All	0.67	0/9215	0.72	4/12513 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	6	PRO	N-CA-CB	5.67	110.10	103.30
1	D	32	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	E	15[A]	LEU	CA-CB-CG	5.28	127.45	115.30
1	E	15[B]	LEU	CA-CB-CG	5.28	127.45	115.30

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	1465	0	1560	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1480	0	1572	18	0
1	C	1488	0	1585	13	0
1	D	1529	0	1646	38	0
1	E	1509	0	1614	34	0
1	F	1506	0	1613	23	0
2	A	5	0	0	0	0
2	B	15	0	0	2	0
2	C	5	0	0	0	0
2	D	20	0	0	3	0
2	E	20	0	0	4	0
2	F	19	0	0	1	0
3	A	132	0	0	4	0
3	B	132	0	0	8	0
3	C	130	0	0	2	0
3	D	220	0	0	10	1
3	E	198	0	0	4	1
3	F	211	0	0	8	0
All	All	10084	0	9590	136	1

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 (136) 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:15[B]:LEU:HD21	1:E:201:LEU:CD1	1.41	1.46
1:E:15[B]:LEU:CD2	1:E:201:LEU:CD1	2.15	1.25
1:F:18[B]:LEU:HD21	3:F:2098:HOH:O	1.06	1.21
1:D:46:LEU:H	1:D:69:THR:CG2	1.66	1.06
1:A:118:GLU:HG3	3:B:2088:HOH:O	1.57	1.05
1:E:15[B]:LEU:CD2	1:E:201:LEU:HD13	1.89	1.01
1:D:46:LEU:H	1:D:69:THR:HG23	1.19	1.01
1:E:18[A]:LEU:HD21	2:E:2014:PO4:O3	1.61	1.00
1:E:15[B]:LEU:HD21	1:E:201:LEU:HD12	0.98	0.96
1:D:162[A]:GLU:HG3	3:D:2164:HOH:O	1.65	0.96
1:E:15[A]:LEU:HD21	1:E:182:LEU:CD1	2.03	0.89
1:E:15[B]:LEU:CD2	1:E:201:LEU:HD12	1.94	0.86
1:B:115:THR:HG22	1:B:119[B]:ARG:HE	1.42	0.82
1:C:8:ALA:O	1:C:12:GLU:HG2	1.81	0.80
1:D:15[B]:LEU:HD11	1:D:182:LEU:HD11	1.62	0.79
1:C:95:GLU:OE1	1:D:95:GLU:OE1	2.04	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:15[B]:LEU:HD12	1:E:198:LYS:HG2	1.69	0.74
1:E:92:LEU:HD22	3:E:2152:HOH:O	1.87	0.74
1:E:20:THR:HG22	1:E:45:THR:OG1	1.88	0.73
1:B:6:PRO:N	3:B:2145:HOH:O	2.22	0.73
1:D:119[B]:ARG:HH22	1:E:119:ARG:HH21	1.36	0.72
1:D:200:LEU:HD12	3:D:2164:HOH:O	1.89	0.72
1:B:95:GLU:OE1	1:F:95:GLU:OE1	2.08	0.71
1:E:15[B]:LEU:HD23	1:E:201:LEU:CD1	2.21	0.71
1:F:198:LYS:HE2	3:F:2194:HOH:O	1.90	0.69
1:E:15[A]:LEU:HD21	1:E:182:LEU:HD11	1.74	0.69
1:D:46:LEU:N	1:D:69:THR:HG23	2.00	0.68
1:E:15[B]:LEU:HD11	1:E:198:LYS:HA	1.74	0.68
1:E:10:LEU:HD21	1:E:16:LEU:HD13	1.75	0.68
1:A:45[B]:THR:HG23	1:A:69:THR:OG1	1.95	0.66
1:D:200:LEU:CD1	3:D:2164:HOH:O	2.43	0.66
1:A:45[B]:THR:HG21	1:A:47:ARG:NE	2.10	0.65
1:C:18:LEU:HD22	1:C:45:THR:CG2	2.26	0.65
1:E:15[B]:LEU:HD23	1:E:201:LEU:HD13	1.75	0.65
1:A:95:GLU:OE1	1:E:95:GLU:OE1	2.15	0.64
1:B:119[B]:ARG:HH22	1:C:119:ARG:CZ	2.11	0.63
1:F:161:LYS:H	1:F:164:HIS:HD2	1.45	0.63
1:C:48:THR:HG23	1:C:50:LYS:H	1.66	0.61
1:D:46:LEU:N	1:D:69:THR:CG2	2.51	0.60
1:E:15[B]:LEU:CD1	1:E:198:LYS:HG2	2.31	0.60
1:E:167:HIS:HB2	3:E:2212:HOH:O	2.03	0.59
1:E:92:LEU:HD23	1:E:119:ARG:HD2	1.84	0.59
1:D:119[B]:ARG:NH2	1:E:119:ARG:HH21	2.01	0.58
1:D:10:LEU:HD21	1:D:16:LEU:HB2	1.85	0.58
1:F:161:LYS:H	1:F:164:HIS:CD2	2.21	0.57
1:D:18[A]:LEU:HD11	1:D:45:THR:CG2	2.34	0.57
1:F:164:HIS:HE1	3:F:2207:HOH:O	1.86	0.56
1:F:116:GLU:HA	1:F:119[A]:ARG:HD3	1.88	0.56
1:B:119[A]:ARG:HG3	3:B:2049:HOH:O	2.05	0.56
1:B:184:GLN:O	1:B:193:LYS:NZ	2.36	0.56
1:F:157[B]:THR:HG21	3:F:2142:HOH:O	2.05	0.56
1:E:111:VAL:HG12	3:F:2149:HOH:O	2.05	0.56
1:E:155:LEU:HD13	3:E:2205:HOH:O	2.04	0.55
1:B:129[B]:LYS:HE3	1:B:131:PHE:CB	2.36	0.55
1:E:55:LEU:HD23	1:E:65[B]:LEU:HD23	1.88	0.55
1:B:110:GLY:HA2	1:B:129[B]:LYS:HB3	1.88	0.55
1:E:21:VAL:HG22	2:E:2004:PO4:O1	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:195:ARG:NH1	3:D:2180:HOH:O	2.39	0.54
1:D:160:ILE:HG21	1:D:177[A]:VAL:CG1	2.38	0.54
1:F:18[B]:LEU:CD2	3:F:2098:HOH:O	1.93	0.53
1:F:10:LEU:HD21	1:F:16:LEU:HD13	1.90	0.53
1:E:47:ARG:HG2	3:F:2107:HOH:O	2.09	0.53
1:F:182:LEU:HD23	1:F:182:LEU:H	1.73	0.53
1:E:15[B]:LEU:CD2	1:E:201:LEU:HD11	2.32	0.51
1:A:129[B]:LYS:HE3	1:A:157:THR:HB	1.91	0.51
1:D:186:ASN:C	1:D:186:ASN:HD22	2.15	0.51
1:D:18[B]:LEU:HG	3:D:2202:HOH:O	2.10	0.51
1:E:182:LEU:HD23	1:E:182:LEU:H	1.75	0.50
1:B:129[B]:LYS:HE3	1:B:131:PHE:HB3	1.93	0.50
1:B:7:LEU:HD22	1:B:10:LEU:HD12	1.92	0.50
3:B:2092:HOH:O	1:C:92:LEU:HB2	2.10	0.50
1:A:41:ALA:HB3	3:A:2083:HOH:O	2.12	0.50
1:D:46:LEU:H	1:D:69:THR:HG21	1.67	0.50
1:E:157[A]:THR:HG22	2:E:2014:PO4:O1	2.12	0.50
1:D:18[A]:LEU:HD11	1:D:45:THR:HG23	1.94	0.50
1:F:15[A]:LEU:HD21	1:F:201:LEU:CD1	2.42	0.49
1:A:45[B]:THR:CG2	1:A:47:ARG:HG3	2.42	0.49
1:C:18:LEU:HD22	1:C:45:THR:HG23	1.94	0.48
1:B:116:GLU:HA	1:B:119[B]:ARG:HD2	1.95	0.48
1:D:160:ILE:HG21	1:D:177[A]:VAL:HG11	1.95	0.48
1:B:129[B]:LYS:NZ	3:B:2143:HOH:O	2.47	0.48
1:D:161:LYS:H	1:D:164:HIS:HD2	1.62	0.47
1:A:110:GLY:HA2	1:A:129[B]:LYS:HB3	1.95	0.47
1:D:157[B]:THR:HG21	3:D:2231:HOH:O	2.14	0.47
1:D:148[B]:VAL:HG23	2:F:1169:PO4:O1	2.13	0.47
1:E:99:LEU:C	1:E:99:LEU:HD13	2.35	0.47
1:D:160:ILE:HD13	1:D:177[A]:VAL:HG13	1.97	0.47
1:F:32:ARG:NH2	3:F:2185:HOH:O	2.46	0.47
1:F:115:THR:HG22	1:F:119[A]:ARG:HD2	1.97	0.47
1:B:88:VAL:HG11	1:B:129[A]:LYS:HE3	1.97	0.46
1:C:7:LEU:HD22	1:C:10:LEU:HD12	1.96	0.46
1:E:161:LYS:HE2	3:E:2149:HOH:O	2.14	0.46
1:B:119[B]:ARG:CZ	2:B:2006:PO4:O1	2.64	0.46
1:A:108:LEU:N	1:A:108:LEU:HD12	2.30	0.46
1:A:10:LEU:HD21	1:A:16:LEU:HD13	1.97	0.46
1:D:144:ALA:O	1:D:147[A]:GLU:HG3	2.15	0.46
1:B:18:LEU:HD21	1:B:45:THR:HG21	1.98	0.46
1:D:69:THR:HG22	2:D:2012:PO4:O4	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:HIS:HB2	3:A:2137:HOH:O	2.16	0.45
1:E:37:GLU:OE1	1:E:195:ARG:HG2	2.17	0.45
1:D:15[B]:LEU:CD1	1:D:182:LEU:HD11	2.38	0.45
1:B:193:LYS:HE2	3:B:2047:HOH:O	2.17	0.45
3:B:2092:HOH:O	1:C:111:VAL:HG12	2.15	0.45
1:E:71[A]:ARG:HB2	1:E:75:GLU:OE1	2.16	0.45
1:F:37:GLU:O	1:F:198:LYS:HD2	2.16	0.45
1:D:101:GLN:HB3	1:D:101:GLN:HE21	1.62	0.44
1:C:48:THR:HG22	3:C:2130:HOH:O	2.16	0.44
1:A:118:GLU:HG3	3:A:2009:HOH:O	2.17	0.43
1:B:119[B]:ARG:NH2	2:B:2006:PO4:O1	2.52	0.43
1:F:95:GLU:OE1	1:F:95:GLU:HA	2.18	0.43
1:F:144:ALA:O	1:F:147:GLU:HG3	2.18	0.43
3:D:2167:HOH:O	1:F:69:THR:HG23	2.19	0.43
1:C:18:LEU:CD2	1:C:45:THR:CG2	2.96	0.42
1:D:145:TYR:HA	1:D:148[B]:VAL:HG12	2.00	0.42
1:D:18[B]:LEU:HD21	3:D:2179:HOH:O	2.19	0.42
1:F:182:LEU:CD2	1:F:182:LEU:H	2.32	0.42
1:A:88:VAL:HG11	1:A:129[A]:LYS:HE3	2.02	0.42
1:D:139:VAL:HG11	1:D:167:HIS:HB3	2.00	0.42
1:D:162[B]:GLU:HG2	3:D:2018:HOH:O	2.18	0.42
1:F:15[A]:LEU:HD21	1:F:201:LEU:HD12	2.02	0.42
1:D:119[B]:ARG:HH22	1:E:119:ARG:NH2	2.11	0.42
1:D:99:LEU:C	1:D:99:LEU:HD13	2.41	0.42
1:B:119[B]:ARG:CZ	3:B:2087:HOH:O	2.67	0.41
1:D:161:LYS:H	1:D:164:HIS:CD2	2.39	0.41
1:D:119[A]:ARG:HG3	3:D:2042:HOH:O	2.19	0.41
1:A:28:LEU:HD22	3:A:2097:HOH:O	2.20	0.41
1:C:44[B]:ILE:HD12	1:C:55:LEU:HD23	2.01	0.41
1:D:69:THR:HB	2:D:2012:PO4:O4	2.21	0.41
1:F:131:PHE:CD1	1:F:132:PRO:HA	2.56	0.41
1:D:69:THR:CG2	2:D:2012:PO4:O4	2.69	0.41
1:F:15[B]:LEU:HD21	1:F:182:LEU:CD1	2.51	0.41
2:E:2010:PO4:O4	2:E:2014:PO4:O1	2.39	0.40
1:A:18:LEU:HD13	1:A:43:GLU:CD	2.41	0.40
1:C:101:GLN:NE2	3:C:2051:HOH:O	2.53	0.40
1:F:166:PRO:HG3	1:F:200:LEU:HD22	2.04	0.40
1:A:36:GLU:OE1	1:A:191:ARG:NH2	2.35	0.40

All (1) 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)
3:D:2230:HOH:O	3:E:2209:HOH:O[3_555]	1.83	0.37

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	196/207 (95%)	192 (98%)	4 (2%)	0	100	100
1	B	198/207 (96%)	193 (98%)	5 (2%)	0	100	100
1	C	199/207 (96%)	194 (98%)	5 (2%)	0	100	100
1	D	207/207 (100%)	205 (99%)	2 (1%)	0	100	100
1	E	203/207 (98%)	199 (98%)	4 (2%)	0	100	100
1	F	202/207 (98%)	199 (98%)	3 (2%)	0	100	100
All	All	1205/1242 (97%)	1182 (98%)	23 (2%)	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	146/153 (95%)	144 (99%)	2 (1%)	74	57
1	B	147/153 (96%)	144 (98%)	3 (2%)	63	41
1	C	150/153 (98%)	148 (99%)	2 (1%)	76	60
1	D	157/153 (103%)	151 (96%)	6 (4%)	40	15

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	152/153 (99%)	147 (97%)	5 (3%)	45	20
1	F	153/153 (100%)	144 (94%)	9 (6%)	24	6
All	All	905/918 (99%)	878 (97%)	27 (3%)	52	24

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	THR
1	A	92	LEU
1	B	14	ARG
1	B	49	GLU
1	B	94	GLU
1	C	18	LEU
1	C	101	GLN
1	D	69	THR
1	D	101	GLN
1	D	147[A]	GLU
1	D	147[B]	GLU
1	D	155	LEU
1	D	186	ASN
1	E	30	LEU
1	E	69	THR
1	E	94	GLU
1	E	137	GLN
1	E	155	LEU
1	F	18[A]	LEU
1	F	18[B]	LEU
1	F	60	LYS
1	F	94	GLU
1	F	119[A]	ARG
1	F	119[B]	ARG
1	F	147	GLU
1	F	148	VAL
1	F	155	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	101	GLN
1	C	137	GLN

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Mol	Chain	Res	Type
1	C	184	GLN
1	D	101	GLN
1	D	164	HIS
1	D	184	GLN
1	D	186	ASN
1	E	137	GLN
1	E	184	GLN
1	F	164	HIS
1	F	184	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

17 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	PO4	A	2005	-	4,4,4	0.92	0	6,6,6	0.33	0
2	PO4	B	2003	-	4,4,4	0.70	0	6,6,6	0.29	0
2	PO4	B	2006	-	4,4,4	0.40	0	6,6,6	0.40	0
2	PO4	B	2015	-	4,4,4	0.56	0	6,6,6	0.27	0
2	PO4	C	2016	-	4,4,4	0.50	0	6,6,6	0.29	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	D	2001	-	4,4,4	0.53	0	6,6,6	0.30	0
2	PO4	D	2002	-	4,4,4	0.34	0	6,6,6	0.28	0
2	PO4	D	2011	-	4,4,4	0.47	0	6,6,6	0.26	0
2	PO4	D	2012	-	4,4,4	0.58	0	6,6,6	0.30	0
2	PO4	E	2004	-	4,4,4	0.33	0	6,6,6	0.26	0
2	PO4	E	2010	-	4,4,4	0.57	0	6,6,6	0.31	0
2	PO4	E	2013	-	4,4,4	0.43	0	6,6,6	0.29	0
2	PO4	E	2014	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	F	1169	1	0,3,4	0.00	-	0,3,6	0.00	-
2	PO4	F	2007	-	4,4,4	0.31	0	6,6,6	0.27	0
2	PO4	F	2008	-	4,4,4	0.25	0	6,6,6	0.39	0
2	PO4	F	2009	-	4,4,4	0.50	0	6,6,6	0.33	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PO4	A	2005	-	-	0/0/0/0	0/0/0/0
2	PO4	B	2003	-	-	0/0/0/0	0/0/0/0
2	PO4	B	2006	-	-	0/0/0/0	0/0/0/0
2	PO4	B	2015	-	-	0/0/0/0	0/0/0/0
2	PO4	C	2016	-	-	0/0/0/0	0/0/0/0
2	PO4	D	2001	-	-	0/0/0/0	0/0/0/0
2	PO4	D	2002	-	-	0/0/0/0	0/0/0/0
2	PO4	D	2011	-	-	0/0/0/0	0/0/0/0
2	PO4	D	2012	-	-	0/0/0/0	0/0/0/0
2	PO4	E	2004	-	-	0/0/0/0	0/0/0/0
2	PO4	E	2010	-	-	0/0/0/0	0/0/0/0
2	PO4	E	2013	-	-	0/0/0/0	0/0/0/0
2	PO4	E	2014	-	-	0/0/0/0	0/0/0/0
2	PO4	F	1169	1	-	0/0/0/0	0/0/0/0
2	PO4	F	2007	-	-	0/0/0/0	0/0/0/0
2	PO4	F	2008	-	-	0/0/0/0	0/0/0/0
2	PO4	F	2009	-	-	0/0/0/0	0/0/0/0

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.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2006	PO4	2	0
2	D	2012	PO4	3	0
2	E	2004	PO4	1	0
2	E	2010	PO4	1	0
2	E	2014	PO4	3	0
2	F	1169	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 will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.