Full wwPDB X-ray Structure Validation Report (i)

Sep 12, 2023 – 05:20 AM EDT

PDB ID	:	4LHP
Title	:	Crystal Structure of Native FG41Malonate Semialdehyde Decarboxylase
Authors	:	Guo, Y.; Serrano, H.; Poelarends, G.J.; Johnson Jr., W.H.; Hackert, M.L.; Whitman, C.P.
Deposited on	:	2013-07-01
Resolution	:	2.02 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

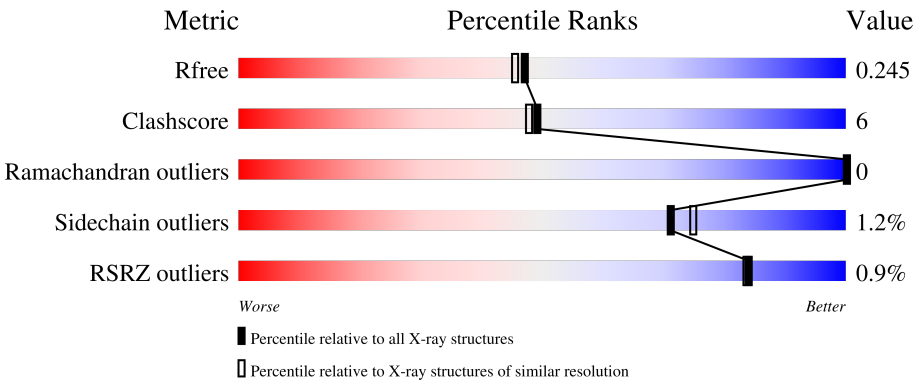
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.02 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	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 >=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% 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	136	90% 5% • •
1	B	136	87% 10% •
1	C	136	86% 10% •
1	D	136	% 85% 11% •
1	E	136	81% 15% •

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Mol	Chain	Length	Quality of chain
1	F	136	85% 10%
1	G	136	% 84% 12%
1	H	136	% 85% 10%
1	I	136	% 81% 15%
1	J	136	4% 79% 17%
1	K	136	% 85% 11%
1	L	136	% 87% 10%

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	F	202	-	-	X	-
2	PO4	G	201	-	-	X	X
2	PO4	H	201	-	-	X	-
2	PO4	J	201	-	-	X	-
2	PO4	K	201	-	-	X	-
2	PO4	L	201	-	-	X	-
3	SO4	E	201	-	-	X	-

2 Entry composition (i)

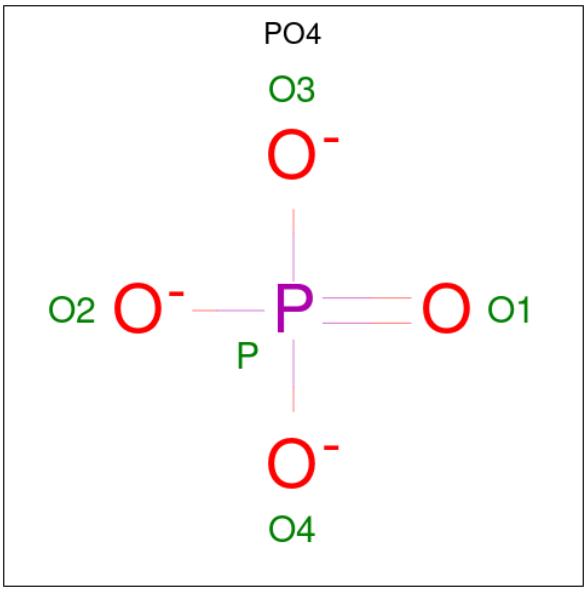
There are 4 unique types of molecules in this entry. The entry contains 13048 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 FG41 Malonate Semialdehyde Decarboxylase.

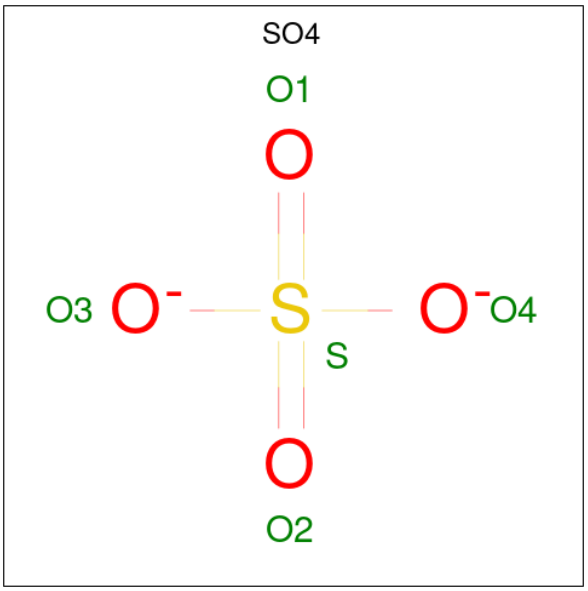
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	131	984	625	175	184	0	0	0
1	B	131	988	627	175	186	0	0	0
1	C	131	988	627	175	186	0	0	0
1	D	131	992	629	175	188	0	0	0
1	E	131	989	628	175	186	0	0	0
1	F	130	989	626	174	189	0	0	0
1	G	131	981	624	174	183	0	0	0
1	H	131	992	629	175	188	0	0	0
1	I	131	969	618	169	182	0	0	0
1	J	131	980	623	174	183	0	0	0
1	K	130	981	622	174	185	0	0	0
1	L	131	972	619	169	184	0	0	0

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	5	4	1	0	0
2	B	1	5	4	1	0	0
2	C	1	5	4	1	0	0
2	D	1	5	4	1	0	0
2	E	1	5	4	1	0	0
2	F	1	5	4	1	0	0
2	G	1	5	4	1	0	0
2	H	1	5	4	1	0	0
2	I	1	5	4	1	0	0
2	J	1	5	4	1	0	0
2	K	1	5	4	1	0	0
2	L	1	5	4	1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total O S 5 4 1	0	0
3	F	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	98	Total O 98 98	0	0
4	B	104	Total O 104 104	0	0
4	C	102	Total O 102 102	0	0
4	D	96	Total O 96 96	0	0
4	E	123	Total O 123 123	0	0
4	F	106	Total O 106 106	0	0
4	G	89	Total O 89 89	0	0
4	H	90	Total O 90 90	0	0
4	I	94	Total O 94 94	0	0
4	J	80	Total O 80 80	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	K	95	95	95	0	0
			Total	O		
4	L	96	96	96	0	0

3 Residue-property plots (i)

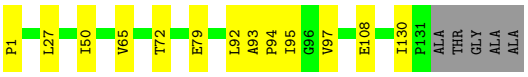
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: FG41 Malonate Semialdehyde Decarboxylase

Chain A:

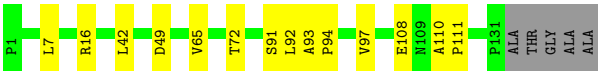
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain B:



• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain C:

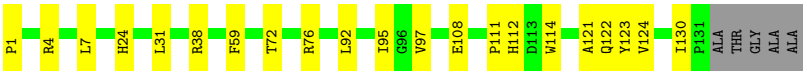


• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain D:

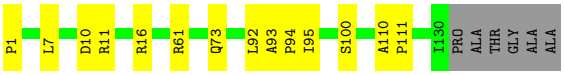
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain E:



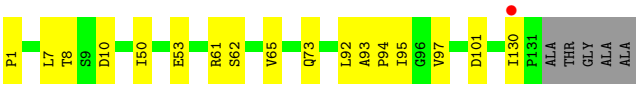
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain F:



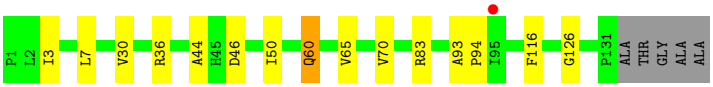
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain G:



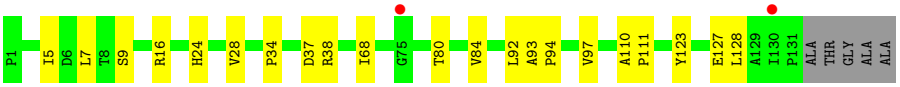
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain H:



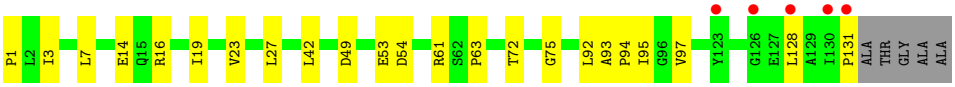
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain I:



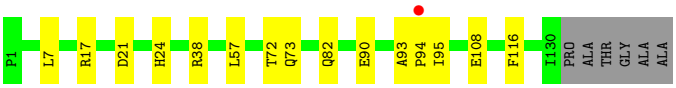
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain J:



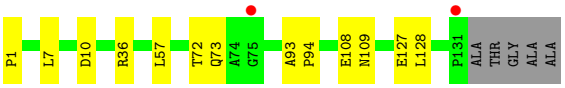
• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain K:



• Molecule 1: FG41 Malonate Semialdehyde Decarboxylase

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	88.95Å 94.69Å 190.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.83 – 2.02 42.02 – 2.02	Depositor EDS
% Data completeness (in resolution range)	85.8 (40.83-2.02) 85.8 (42.02-2.02)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$< I/\sigma(I) >$ [1]	1.87 (at 2.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.194 , 0.241 0.201 , 0.245	Depositor DCC
R_{free} test set	4565 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	21.2	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 54.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	13048	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 53.65 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.0903e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	A	0.22	0/1004	0.45	0/1370
1	B	0.24	0/1008	0.46	0/1375
1	C	0.23	0/1008	0.45	0/1375
1	D	0.23	0/1012	0.46	0/1380
1	E	0.23	0/1009	0.48	0/1376
1	F	0.23	0/1008	0.45	0/1373
1	G	0.22	0/1001	0.44	0/1366
1	H	0.21	0/1012	0.44	0/1380
1	I	0.25	0/989	0.48	0/1352
1	J	0.23	0/1000	0.44	0/1365
1	K	0.24	0/1000	0.47	0/1363
1	L	0.22	0/992	0.45	0/1356
All	All	0.23	0/12043	0.46	0/16431

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	984	0	975	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	988	0	979	10	0
1	C	988	0	979	11	0
1	D	992	0	983	8	0
1	E	989	0	981	20	0
1	F	989	0	980	9	0
1	G	981	0	971	13	0
1	H	992	0	983	13	0
1	I	969	0	951	11	0
1	J	980	0	969	19	0
1	K	981	0	972	14	0
1	L	972	0	953	11	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	1	0
2	F	5	0	0	2	0
2	G	5	0	0	2	0
2	H	5	0	0	2	0
2	I	5	0	0	0	0
2	J	5	0	0	4	0
2	K	5	0	0	2	0
2	L	5	0	0	3	0
3	E	5	0	0	10	0
3	F	5	0	0	0	0
4	A	98	0	0	2	0
4	B	104	0	0	2	0
4	C	102	0	0	5	0
4	D	96	0	0	1	0
4	E	123	0	0	1	0
4	F	106	0	0	1	0
4	G	89	0	0	2	0
4	H	90	0	0	8	0
4	I	94	0	0	1	1
4	J	80	0	0	8	0
4	K	95	0	0	6	0
4	L	96	0	0	2	1
All	All	13048	0	11676	140	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 6.

All (140) 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:C:16:ARG:HD2	4:C:400:HOH:O	1.16	1.29
1:F:1:PRO:N	2:F:202:PO4:O2	1.87	1.08
1:F:73:GLN:NE2	2:F:202:PO4:O1	2.04	0.90
1:E:123:TYR:N	3:E:201:SO4:O1	2.07	0.88
1:H:44:ALA:O	4:H:390:HOH:O	1.96	0.84
1:C:65:VAL:O	4:C:352:HOH:O	1.95	0.83
1:J:53:GLU:O	1:J:61:ARG:NH2	2.14	0.81
2:H:201:PO4:O3	4:H:381:HOH:O	1.98	0.80
1:D:5:ILE:HB	1:D:42:LEU:HD23	1.68	0.76
1:H:60:GLN:OE1	4:H:364:HOH:O	2.04	0.76
1:E:124:VAL:HG23	3:E:201:SO4:O4	1.86	0.76
1:J:14:GLU:O	4:J:378:HOH:O	2.01	0.76
1:G:53:GLU:O	1:G:61:ARG:NH1	2.21	0.73
1:E:121:ALA:C	3:E:201:SO4:O1	2.28	0.71
1:J:63:PRO:O	4:J:330:HOH:O	2.08	0.71
1:C:49:ASP:OD2	4:C:356:HOH:O	2.12	0.67
1:C:91:SER:O	4:C:370:HOH:O	2.13	0.67
1:F:92:LEU:HD23	1:F:95:ILE:HD11	1.76	0.67
1:C:49:ASP:OD2	4:C:387:HOH:O	2.12	0.66
1:B:50:ILE:HG21	1:B:65:VAL:CG2	2.26	0.66
1:J:49:ASP:OD2	4:J:370:HOH:O	2.12	0.66
1:B:79:GLU:OE2	4:B:360:HOH:O	2.13	0.66
1:E:121:ALA:HB1	3:E:201:SO4:O3	1.96	0.65
1:E:122:GLN:N	3:E:201:SO4:O1	2.31	0.63
1:K:82:GLN:O	4:K:388:HOH:O	2.16	0.61
1:J:93:ALA:HB3	1:J:94:PRO:HD3	1.84	0.60
1:K:17:ARG:NH1	4:K:374:HOH:O	2.34	0.60
1:K:82:GLN:C	4:K:388:HOH:O	2.40	0.59
1:I:34:PRO:HG2	1:I:37:ASP:HB2	1.85	0.58
1:H:36:ARG:HG3	4:H:349:HOH:O	2.03	0.58
1:H:46:ASP:HB2	4:H:370:HOH:O	2.03	0.58
1:E:1:PRO:N	4:E:380:HOH:O	2.22	0.57
1:J:61:ARG:NH1	1:K:116:PHE:O	2.37	0.57
1:I:123:TYR:CE2	1:I:128:LEU:HD23	2.40	0.57
1:H:30:VAL:HG13	1:H:83:ARG:NH1	2.20	0.56
1:J:16:ARG:NH1	4:J:345:HOH:O	2.37	0.56
1:K:73:GLN:NE2	2:K:201:PO4:O2	2.40	0.55
1:G:62:SER:OG	1:G:101:ASP:OD2	2.15	0.55
1:E:114:TRP:O	3:E:201:SO4:O3	2.24	0.55
1:J:75:GLY:HA3	4:J:354:HOH:O	2.07	0.55
1:K:21:ASP:HB2	4:K:313:HOH:O	2.07	0.55
1:G:50:ILE:HG21	1:G:65:VAL:CG2	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:GLU:OE2	4:D:364:HOH:O	2.18	0.54
1:L:73:GLN:NE2	2:L:201:PO4:O2	2.39	0.54
1:J:72:THR:OG1	2:J:201:PO4:O3	2.22	0.54
1:E:112:HIS:HA	3:E:201:SO4:O4	2.07	0.54
1:I:24:HIS:CE1	1:I:38:ARG:HB2	2.43	0.54
1:G:130:ILE:O	4:G:387:HOH:O	2.18	0.54
1:I:127:GLU:OE1	4:I:332:HOH:O	2.18	0.54
1:H:3:ILE:HD12	1:H:70:VAL:HG22	1.90	0.53
1:E:123:TYR:HB2	3:E:201:SO4:O4	2.10	0.52
4:K:319:HOH:O	1:L:109:ASN:HB3	2.10	0.52
1:E:31:LEU:HD22	1:E:76:ARG:HD2	1.91	0.52
1:D:121:ALA:O	1:D:124:VAL:HG22	2.10	0.52
1:G:1:PRO:H3	2:G:201:PO4:P	2.33	0.52
1:B:50:ILE:CG2	1:B:65:VAL:HG21	2.40	0.51
1:K:93:ALA:HB3	1:K:94:PRO:HD3	1.92	0.50
1:K:73:GLN:N	2:K:201:PO4:O4	2.45	0.50
1:K:93:ALA:O	1:K:95:ILE:O	2.30	0.50
1:E:72:THR:OG1	2:E:202:PO4:O2	2.19	0.50
1:D:93:ALA:N	1:D:94:PRO:HD2	2.26	0.49
1:E:111:PRO:O	3:E:201:SO4:O4	2.31	0.49
1:B:50:ILE:HG21	1:B:65:VAL:HG21	1.94	0.49
1:J:54:ASP:HB3	4:J:317:HOH:O	2.13	0.48
1:L:93:ALA:N	1:L:94:PRO:HD2	2.28	0.48
1:I:93:ALA:N	1:I:94:PRO:HD2	2.29	0.48
1:H:93:ALA:HB3	1:H:94:PRO:CD	2.45	0.47
1:A:95:ILE:HD11	4:A:326:HOH:O	2.14	0.47
1:L:72:THR:O	1:L:108:GLU:HA	2.15	0.47
4:K:383:HOH:O	1:L:36:ARG:CB	2.63	0.47
1:D:92:LEU:HB3	1:D:97:VAL:HB	1.97	0.47
1:E:59:PHE:HB2	4:L:392:HOH:O	2.14	0.47
1:B:72:THR:O	1:B:108:GLU:HA	2.15	0.47
1:K:57:LEU:HD21	1:L:128:LEU:HG	1.97	0.47
1:J:131:PRO:HB3	4:J:363:HOH:O	2.14	0.46
1:C:93:ALA:HB3	1:C:94:PRO:HD3	1.97	0.46
1:G:95:ILE:HG23	4:G:385:HOH:O	2.14	0.46
1:C:92:LEU:HB3	1:C:97:VAL:HB	1.97	0.46
1:K:90:GLU:O	1:K:94:PRO:HD3	2.16	0.46
1:L:72:THR:OG1	2:L:201:PO4:O1	2.28	0.46
1:C:42:LEU:N	1:C:42:LEU:CD1	2.78	0.46
1:G:73:GLN:NE2	2:G:201:PO4:O3	2.46	0.46
1:K:24:HIS:CE1	1:K:38:ARG:HB2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:1:PRO:HA	2:L:201:PO4:O3	2.16	0.46
1:J:19:ILE:O	1:J:23:VAL:HG23	2.16	0.46
1:F:110:ALA:HB1	1:F:111:PRO:CD	2.45	0.46
1:G:61:ARG:NH2	1:H:116:PHE:O	2.48	0.45
1:I:80:THR:O	1:I:84:VAL:HG23	2.15	0.45
1:F:16:ARG:NH2	4:F:334:HOH:O	2.50	0.45
1:F:93:ALA:N	1:F:94:PRO:CD	2.80	0.45
1:J:95:ILE:HD11	4:J:318:HOH:O	2.16	0.45
1:D:6:ASP:OD2	1:E:4:ARG:NH2	2.49	0.45
1:A:92:LEU:HB3	1:A:97:VAL:HB	1.99	0.45
1:I:110:ALA:HB1	1:I:111:PRO:HD2	1.99	0.45
1:B:95:ILE:HG22	1:B:95:ILE:O	2.16	0.45
1:J:128:LEU:HB2	1:L:57:LEU:HD21	1.98	0.45
1:E:95:ILE:HG22	1:E:95:ILE:O	2.17	0.44
1:E:72:THR:O	1:E:108:GLU:HA	2.17	0.44
2:H:201:PO4:P	4:H:381:HOH:O	2.69	0.44
1:H:93:ALA:N	1:H:94:PRO:HD2	2.32	0.44
1:F:10:ASP:OD2	1:F:11:ARG:N	2.51	0.44
1:F:61:ARG:NH1	1:F:100:SER:O	2.51	0.43
1:G:50:ILE:CG2	1:G:65:VAL:CG2	2.96	0.43
1:G:93:ALA:HB3	1:G:94:PRO:HD3	2.00	0.43
1:C:110:ALA:HB1	1:C:111:PRO:HD2	2.00	0.43
1:H:50:ILE:HG21	1:H:65:VAL:CG2	2.47	0.43
1:J:3:ILE:HD11	1:J:27:LEU:CD1	2.49	0.43
1:J:92:LEU:HB3	1:J:97:VAL:HB	1.99	0.43
1:A:93:ALA:N	1:A:94:PRO:HD2	2.33	0.43
1:E:111:PRO:HB3	1:E:130:ILE:CD1	2.49	0.43
1:I:92:LEU:HB3	1:I:97:VAL:HB	2.01	0.43
1:K:72:THR:O	1:K:108:GLU:HA	2.17	0.43
1:B:93:ALA:HB3	1:B:94:PRO:HD3	1.99	0.43
1:E:92:LEU:HB3	1:E:97:VAL:HB	2.01	0.43
1:D:110:ALA:HB1	1:D:111:PRO:HD2	2.00	0.43
1:I:24:HIS:O	1:I:28:VAL:HG23	2.19	0.42
1:C:93:ALA:N	1:C:94:PRO:HD2	2.34	0.42
1:D:95:ILE:O	1:D:95:ILE:HG22	2.19	0.42
1:H:93:ALA:HB3	1:H:94:PRO:HD3	2.01	0.42
1:A:7:LEU:HD12	1:A:44:ALA:HA	2.02	0.42
1:E:123:TYR:HB2	3:E:201:SO4:S	2.60	0.42
1:G:92:LEU:HB3	1:G:97:VAL:HB	2.02	0.42
1:E:24:HIS:CE1	1:E:38:ARG:HB2	2.55	0.42
1:H:126:GLY:HA2	4:H:308:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:72:THR:HA	2:J:201:PO4:O3	2.20	0.41
1:I:5:ILE:HG12	1:I:68:ILE:HD12	2.02	0.41
1:J:1:PRO:HB3	2:J:201:PO4:O3	2.21	0.41
1:B:1:PRO:N	4:B:391:HOH:O	2.53	0.41
1:B:92:LEU:HB3	1:B:97:VAL:HB	2.02	0.41
1:L:10:ASP:OD2	4:L:394:HOH:O	2.22	0.41
1:G:8:THR:OG1	1:G:10:ASP:OD2	2.20	0.41
1:K:57:LEU:HD22	1:L:127:GLU:HG2	2.03	0.41
1:J:1:PRO:CA	2:J:201:PO4:O3	2.69	0.41
1:H:83:ARG:HD3	4:H:339:HOH:O	2.21	0.40
1:I:9:SER:HA	1:I:16:ARG:NH2	2.36	0.40
1:C:72:THR:O	1:C:108:GLU:HA	2.21	0.40
1:F:93:ALA:HB3	1:F:94:PRO:HD3	2.03	0.40
1:G:50:ILE:HG21	1:G:65:VAL:HG23	2.04	0.40
1:A:91:SER:HB3	4:A:385:HOH:O	2.22	0.40
1:B:130:ILE:O	1:B:130:ILE:HG13	2.20	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)
4:I:371:HOH:O	4:L:375:HOH:O[1_655]	1.89	0.31

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	129/136 (95%)	127 (98%)	2 (2%)	0	100	100
1	B	129/136 (95%)	126 (98%)	3 (2%)	0	100	100
1	C	129/136 (95%)	123 (95%)	6 (5%)	0	100	100
1	D	129/136 (95%)	125 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	129/136 (95%)	125 (97%)	4 (3%)	0	100	100
1	F	128/136 (94%)	124 (97%)	4 (3%)	0	100	100
1	G	129/136 (95%)	124 (96%)	5 (4%)	0	100	100
1	H	129/136 (95%)	125 (97%)	4 (3%)	0	100	100
1	I	129/136 (95%)	126 (98%)	3 (2%)	0	100	100
1	J	129/136 (95%)	128 (99%)	1 (1%)	0	100	100
1	K	128/136 (94%)	124 (97%)	4 (3%)	0	100	100
1	L	129/136 (95%)	127 (98%)	2 (2%)	0	100	100
All	All	1546/1632 (95%)	1504 (97%)	42 (3%)	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	101/106 (95%)	100 (99%)	1 (1%)	76	80
1	B	102/106 (96%)	101 (99%)	1 (1%)	76	80
1	C	102/106 (96%)	101 (99%)	1 (1%)	76	80
1	D	103/106 (97%)	101 (98%)	2 (2%)	57	59
1	E	102/106 (96%)	101 (99%)	1 (1%)	76	80
1	F	103/106 (97%)	102 (99%)	1 (1%)	76	80
1	G	100/106 (94%)	99 (99%)	1 (1%)	76	80
1	H	103/106 (97%)	101 (98%)	2 (2%)	57	59
1	I	98/106 (92%)	97 (99%)	1 (1%)	76	80
1	J	100/106 (94%)	98 (98%)	2 (2%)	55	57
1	K	101/106 (95%)	100 (99%)	1 (1%)	76	80
1	L	99/106 (93%)	98 (99%)	1 (1%)	76	80
All	All	1214/1272 (95%)	1199 (99%)	15 (1%)	71	75

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	B	27	LEU
1	C	7	LEU
1	D	7	LEU
1	D	46	ASP
1	E	7	LEU
1	F	7	LEU
1	G	7	LEU
1	H	7	LEU
1	H	60	GLN
1	I	7	LEU
1	J	7	LEU
1	J	42	LEU
1	K	7	LEU
1	L	7	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	73	GLN
1	H	60	GLN
1	K	73	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)

14 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	G	201	-	4,4,4	0.98	0	6,6,6	0.39	0
2	PO4	J	201	-	4,4,4	0.98	0	6,6,6	0.39	0
3	SO4	E	201	-	4,4,4	0.15	0	6,6,6	0.22	0
2	PO4	A	201	-	4,4,4	0.89	0	6,6,6	0.45	0
3	SO4	F	201	-	4,4,4	0.14	0	6,6,6	0.10	0
2	PO4	C	201	-	4,4,4	0.92	0	6,6,6	0.43	0
2	PO4	F	202	-	4,4,4	0.93	0	6,6,6	0.51	0
2	PO4	K	201	-	4,4,4	0.90	0	6,6,6	0.44	0
2	PO4	B	201	-	4,4,4	0.89	0	6,6,6	0.50	0
2	PO4	D	201	-	4,4,4	0.92	0	6,6,6	0.46	0
2	PO4	E	202	-	4,4,4	0.89	0	6,6,6	0.46	0
2	PO4	I	201	-	4,4,4	0.89	0	6,6,6	0.40	0
2	PO4	L	201	-	4,4,4	0.94	0	6,6,6	0.36	0
2	PO4	H	201	-	4,4,4	0.92	0	6,6,6	0.43	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.

8 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	201	PO4	2	0
2	J	201	PO4	4	0
3	E	201	SO4	10	0
2	F	202	PO4	2	0
2	K	201	PO4	2	0
2	E	202	PO4	1	0
2	L	201	PO4	3	0
2	H	201	PO4	2	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, 95th 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^2)	Q<0.9
1	A	131/136 (96%)	-0.19	0	100	100	13, 23, 32, 46	0
1	B	131/136 (96%)	-0.21	0	100	100	10, 17, 27, 38	0
1	C	131/136 (96%)	-0.20	0	100	100	10, 16, 23, 32	0
1	D	131/136 (96%)	-0.19	2 (1%)	73	73	11, 21, 30, 46	0
1	E	131/136 (96%)	-0.23	0	100	100	10, 17, 27, 46	0
1	F	130/136 (95%)	-0.27	0	100	100	11, 17, 24, 34	0
1	G	131/136 (96%)	-0.05	1 (0%)	86	85	11, 19, 30, 36	0
1	H	131/136 (96%)	-0.06	1 (0%)	86	85	13, 21, 32, 48	0
1	I	131/136 (96%)	-0.21	2 (1%)	73	73	10, 17, 28, 44	0
1	J	131/136 (96%)	0.02	5 (3%)	40	40	12, 21, 29, 37	0
1	K	130/136 (95%)	0.04	1 (0%)	86	85	14, 22, 32, 36	0
1	L	131/136 (96%)	-0.30	2 (1%)	73	73	10, 18, 26, 42	0
All	All	1570/1632 (96%)	-0.16	14 (0%)	84	83	10, 19, 30, 48	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	130	ILE	4.3
1	K	94	PRO	3.7
1	I	75	GLY	3.7
1	L	75	GLY	3.0
1	H	95	ILE	2.4
1	L	131	PRO	2.2
1	J	123	TYR	2.2
1	I	130	ILE	2.1
1	J	126	GLY	2.1
1	D	12	SER	2.1
1	J	131	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	130	ILE	2.1
1	J	128	LEU	2.1
1	D	131	PRO	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th 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^2)	Q<0.9
2	PO4	G	201	5/5	0.66	0.41	33,49,53,63	0
2	PO4	E	202	5/5	0.73	0.35	38,47,50,54	0
2	PO4	C	201	5/5	0.76	0.32	26,44,52,60	0
2	PO4	F	202	5/5	0.82	0.29	43,46,47,52	0
2	PO4	J	201	5/5	0.82	0.34	40,42,48,51	0
2	PO4	L	201	5/5	0.84	0.18	45,46,49,57	0
2	PO4	K	201	5/5	0.86	0.26	44,48,53,62	0
2	PO4	B	201	5/5	0.86	0.36	39,41,54,56	0
3	SO4	E	201	5/5	0.88	0.77	42,48,52,56	0
2	PO4	D	201	5/5	0.89	0.36	41,49,52,59	0
2	PO4	H	201	5/5	0.90	0.25	45,47,54,59	0
2	PO4	A	201	5/5	0.92	0.30	42,48,53,60	0
2	PO4	I	201	5/5	0.92	0.36	39,41,50,52	0
3	SO4	F	201	5/5	0.94	0.17	36,37,48,58	0

6.5 Other polymers (i)

There are no such residues in this entry.