



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:12 AM GMT

PDB ID : 3L6O  
Title : Crystal Structure of Phosphate bound apo Glyceraldehyde-3-phosphate dehydrogenase 1 from MRSA252 at 2.2 Angstrom resolution  
Authors : Mukherjee, S.; Dutta, D.; Saha, B.; Das, A.K.  
Deposited on : 2009-12-23  
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](mailto: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.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

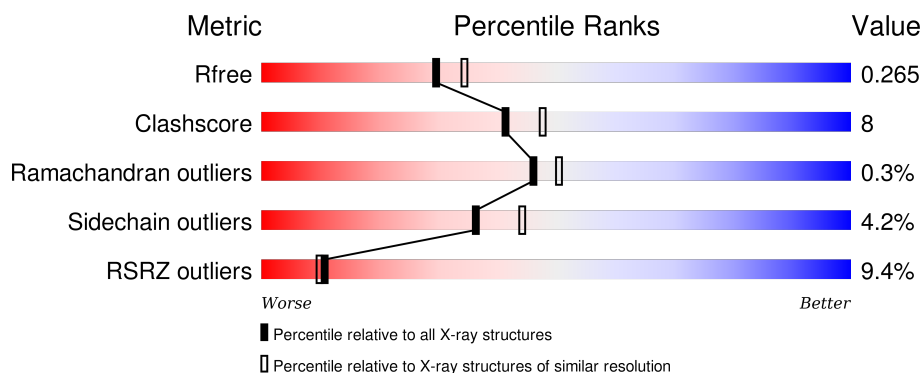
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (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  $\geq 3$ , 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.

Mol	Chain	Length	Quality of chain
1	O	336	<div> <div>13%</div> <div>78%</div> <div>20%</div> <div>..</div> </div>
1	P	336	<div> <div>8%</div> <div>79%</div> <div>19%</div> <div>..</div> </div>
1	Q	336	<div> <div>10%</div> <div>81%</div> <div>16%</div> <div>..</div> </div>
1	R	336	<div> <div>7%</div> <div>81%</div> <div>18%</div> <div>..</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10272 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 Glyceraldehyde-3-phosphate dehydrogenase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	334	Total	C	N	O	S	0	0	0
			2531	1576	435	510	10			
1	P	334	Total	C	N	O	S	0	0	0
			2531	1576	435	510	10			
1	R	334	Total	C	N	O	S	0	0	0
			2531	1576	435	510	10			
1	O	333	Total	C	N	O	S	0	0	0
			2523	1570	434	509	10			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	Q	1	Total	O	P	0	0
			5	4	1		
2	P	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	R	1	Total	O	P	0	0
			5	4	1		
2	O	1	Total	O	P	0	0
			5	4	1		

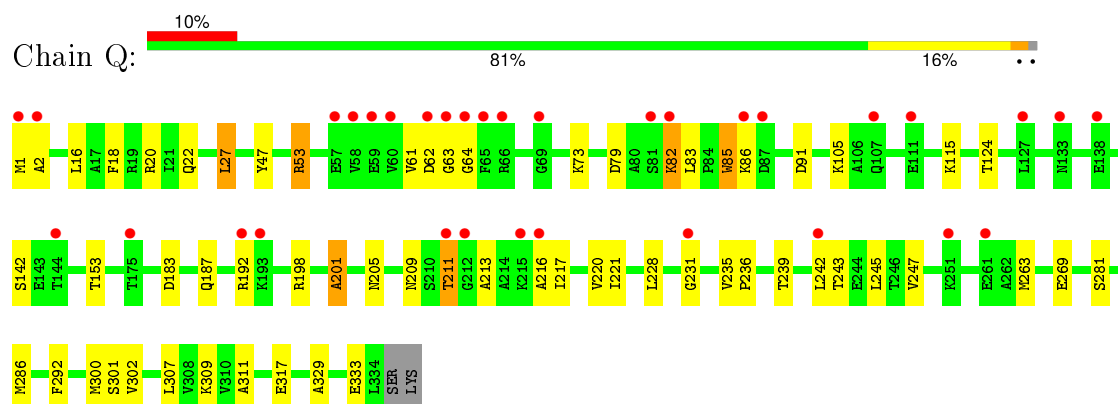
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	Q	33	Total	O	0	0
			33	33		
3	P	39	Total	O	0	0
			39	39		
3	R	35	Total	O	0	0
			35	35		
3	O	29	Total	O	0	0
			29	29		

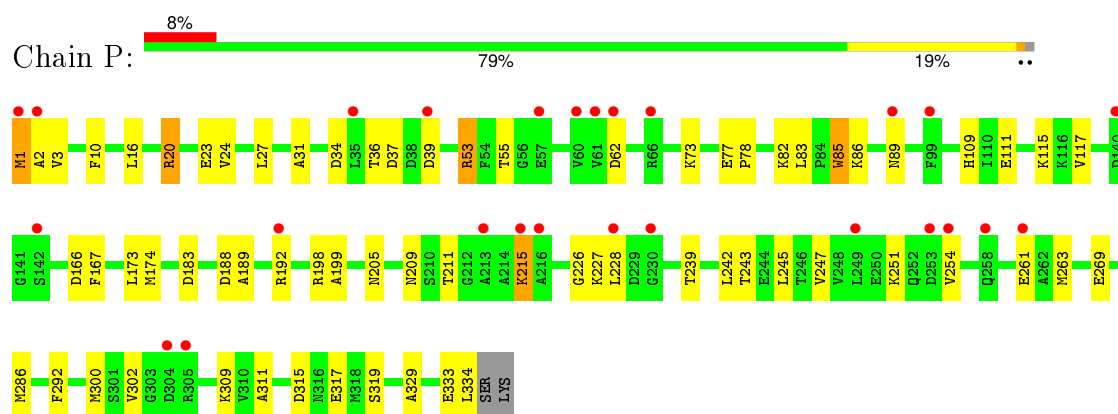
### 3 Residue-property plots

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.

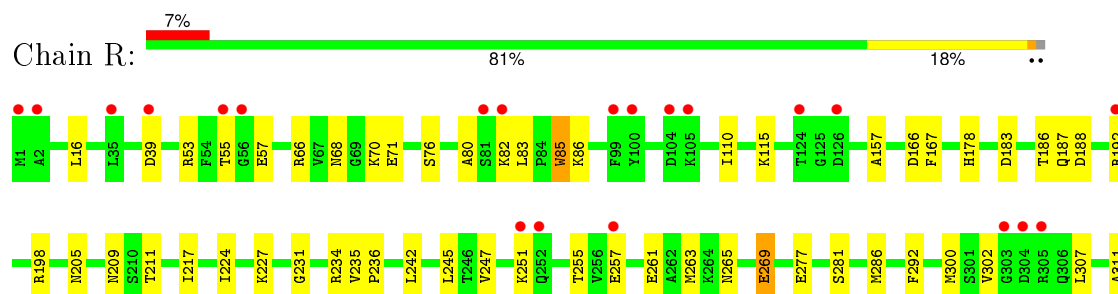
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1

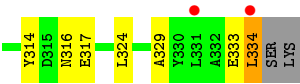


- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1

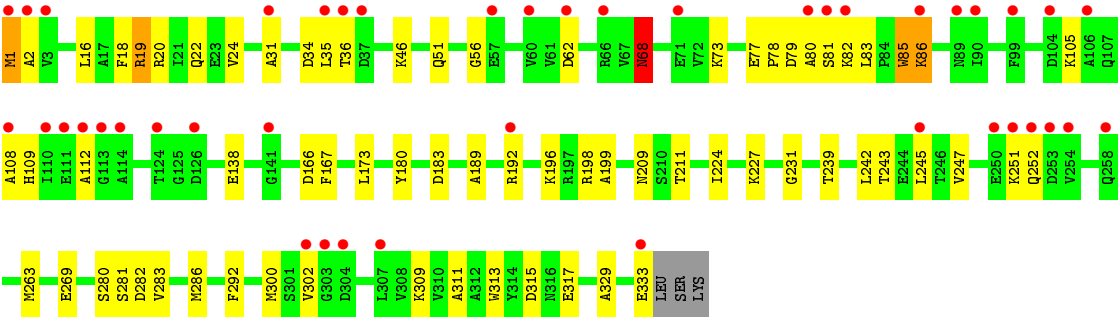
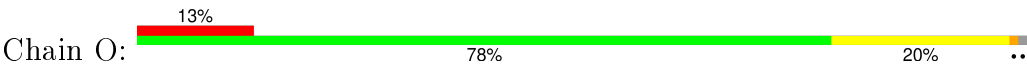


- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1





● Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.95Å 93.68Å 89.05Å 90.00° 106.84° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 33.34 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-2.20) 99.7 (33.34-2.20)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.86 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0095	Depositor
R, $R_{free}$	0.216 , 0.258 0.227 , 0.265	Depositor DCC
$R_{free}$ test set	2706 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	42.4	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 53.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 53347 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10272	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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	O	0.66	0/2557	0.70	1/3464 (0.0%)
1	P	0.69	0/2565	0.74	1/3475 (0.0%)
1	Q	0.70	0/2565	0.76	1/3475 (0.0%)
1	R	0.69	0/2565	0.73	0/3475
All	All	0.69	0/10252	0.73	3/13889 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	27	LEU	CA-CB-CG	6.49	130.22	115.30
1	P	188	ASP	CB-CG-OD1	5.70	123.43	118.30
1	O	19	ARG	NE-CZ-NH2	-5.04	117.78	120.30

There are no chirality outliers.

There are no planarity outliers.

### 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	O	2523	0	2501	60	1
1	P	2531	0	2512	39	3
1	Q	2531	0	2512	47	5
1	R	2531	0	2512	41	4

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	O	5	0	0	0	0
2	P	5	0	0	0	0
2	Q	5	0	0	0	0
2	R	5	0	0	0	0
3	O	29	0	0	15	0
3	P	39	0	0	1	0
3	Q	33	0	0	9	0
3	R	35	0	0	8	0
All	All	10272	0	10037	171	8

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 8.

All (171) 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:Q:187:GLN:HA	3:Q:366:HOH:O	1.25	1.24
1:O:252:GLN:HB2	3:O:353:HOH:O	1.46	1.15
1:O:105:LYS:HE3	3:O:360:HOH:O	1.53	1.07
1:R:324:LEU:HD22	3:R:370:HOH:O	1.53	1.06
1:O:333:GLU:HB3	3:O:361:HOH:O	1.60	1.00
1:Q:53:ARG:HG3	1:Q:53:ARG:HH11	1.35	0.90
1:R:277:GLU:HG3	3:R:361:HOH:O	1.72	0.89
1:O:77:GLU:HG3	1:O:79:ASP:O	1.73	0.88
1:Q:79:ASP:OD1	1:Q:82:LYS:HD3	1.78	0.83
1:O:112:ALA:HB2	3:O:355:HOH:O	1.78	0.83
1:P:53:ARG:HG2	3:P:359:HOH:O	1.78	0.81
1:R:324:LEU:HB2	3:R:370:HOH:O	1.85	0.76
1:Q:79:ASP:CG	1:Q:82:LYS:HD3	2.05	0.76
1:Q:142:SER:HA	3:Q:369:HOH:O	1.86	0.75
1:R:70:LYS:HD2	3:R:364:HOH:O	1.88	0.74
1:Q:105:LYS:HD3	3:Q:361:HOH:O	1.91	0.70
1:Q:198:ARG:HE	1:Q:209:ASN:ND2	1.90	0.69
1:P:3:VAL:HB	1:P:27:LEU:HD23	1.75	0.69
1:P:173:LEU:HD13	1:O:309:LYS:HB2	1.75	0.69
1:O:81:SER:C	3:O:355:HOH:O	2.31	0.68
1:Q:83:LEU:HD13	1:Q:85:TRP:CZ2	2.29	0.67
1:R:198:ARG:HE	1:R:209:ASN:ND2	1.93	0.66
1:P:198:ARG:HE	1:P:209:ASN:ND2	1.92	0.66
1:O:329:ALA:O	1:O:333:GLU:HG3	1.95	0.66
1:R:183:ASP:HB2	1:R:192:ARG:HH22	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:329:ALA:O	1:Q:333:GLU:HG3	1.95	0.65
1:O:16:LEU:HD21	1:O:51:GLN:HG3	1.80	0.64
1:Q:2:ALA:HB3	3:Q:359:HOH:O	1.98	0.64
1:O:77:GLU:OE2	1:O:78:PRO:HD2	1.97	0.64
1:O:198:ARG:HE	1:O:209:ASN:ND2	1.96	0.63
1:O:56:GLY:HA3	1:O:68:ASN:OD1	1.98	0.63
1:O:300:MET:CE	1:O:302:VAL:HG23	2.28	0.63
1:P:183:ASP:HB2	1:P:192:ARG:HH22	1.64	0.62
1:P:1:MET:SD	1:P:2:ALA:N	2.72	0.62
1:O:1:MET:O	1:O:2:ALA:HB3	2.00	0.62
1:Q:307:LEU:HD21	1:R:307:LEU:HD21	1.80	0.62
1:Q:236:PRO:HB2	1:R:236:PRO:HB2	1.82	0.62
1:R:39:ASP:OD2	1:O:196:LYS:NZ	2.27	0.61
1:P:77:GLU:OE2	1:P:78:PRO:HD2	2.00	0.61
1:P:205:ASN:HD22	1:O:281:SER:H	1.49	0.60
1:Q:183:ASP:HB2	1:Q:192:ARG:HH22	1.65	0.60
1:P:83:LEU:HD13	1:P:85:TRP:CZ2	2.37	0.60
1:O:183:ASP:HB2	1:O:192:ARG:HH22	1.65	0.60
1:Q:211:THR:HG22	1:Q:231:GLY:HA2	1.83	0.59
1:O:300:MET:HE2	1:O:302:VAL:HG23	1.84	0.59
1:O:138:GLU:CG	3:O:364:HOH:O	2.50	0.59
1:Q:79:ASP:O	1:Q:82:LYS:HB2	2.03	0.58
1:O:112:ALA:CB	3:O:355:HOH:O	2.45	0.58
1:O:83:LEU:HD23	3:O:362:HOH:O	2.02	0.58
1:Q:300:MET:CE	1:Q:302:VAL:HG23	2.34	0.58
1:Q:198:ARG:HE	1:Q:209:ASN:HD21	1.51	0.58
1:O:211:THR:HG22	1:O:231:GLY:HA2	1.85	0.57
1:P:242:LEU:HD11	1:P:311:ALA:HB1	1.86	0.57
1:R:53:ARG:NH1	3:R:338:HOH:O	2.37	0.56
1:P:227:LYS:HA	1:O:300:MET:HE1	1.87	0.56
1:R:157:ALA:HA	1:R:217:ILE:HD11	1.87	0.56
1:R:257:GLU:HB3	3:R:359:HOH:O	2.06	0.56
1:O:198:ARG:HE	1:O:209:ASN:HD21	1.52	0.55
1:O:35:LEU:HD12	3:O:365:HOH:O	2.06	0.55
1:R:198:ARG:HE	1:R:209:ASN:HD21	1.54	0.55
1:R:329:ALA:O	1:R:333:GLU:HG3	2.06	0.55
1:R:66:ARG:NH1	1:R:71:GLU:HB2	2.21	0.55
1:O:263:MET:HG3	1:O:292:PHE:CE1	2.42	0.54
1:O:46:LYS:CG	3:O:349:HOH:O	2.55	0.54
1:R:263:MET:HG3	1:R:292:PHE:CE1	2.42	0.54
1:P:198:ARG:HE	1:P:209:ASN:HD21	1.54	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:300:MET:HE3	1:P:302:VAL:HG23	1.89	0.54
1:O:86:LYS:HG3	3:O:358:HOH:O	2.07	0.54
1:Q:201:ALA:HB2	3:Q:366:HOH:O	2.08	0.53
1:Q:153:THR:OG1	1:Q:213:ALA:HA	2.08	0.53
1:Q:115:LYS:HE2	3:Q:364:HOH:O	2.07	0.53
1:R:57:GLU:H	1:R:68:ASN:HD21	1.57	0.53
1:R:16:LEU:CD1	1:R:317:GLU:HB3	2.40	0.52
1:Q:183:ASP:HB2	1:Q:192:ARG:NH2	2.24	0.52
1:P:329:ALA:O	1:P:333:GLU:HG3	2.09	0.52
1:O:20:ARG:O	1:O:24:VAL:HG23	2.09	0.52
1:Q:115:LYS:CE	3:Q:364:HOH:O	2.57	0.52
1:R:300:MET:CE	1:R:302:VAL:HG23	2.40	0.52
1:Q:53:ARG:NH1	1:Q:53:ARG:HG3	2.12	0.52
1:R:183:ASP:HB2	1:R:192:ARG:NH2	2.24	0.52
1:Q:281:SER:H	1:R:205:ASN:HD22	1.58	0.51
1:P:183:ASP:OD1	1:P:198:ARG:NH1	2.43	0.51
1:Q:235:VAL:HG11	1:R:235:VAL:HG11	1.93	0.51
1:Q:16:LEU:CD1	1:Q:317:GLU:HB3	2.40	0.51
1:R:211:THR:HG22	1:R:231:GLY:HA2	1.93	0.50
1:P:205:ASN:HD21	1:O:281:SER:CB	2.24	0.50
1:R:316:ASN:ND2	1:R:316:ASN:H	2.10	0.50
1:Q:243:THR:O	1:Q:311:ALA:HA	2.11	0.50
1:O:46:LYS:HG2	3:O:349:HOH:O	2.11	0.50
1:R:300:MET:HE3	1:R:302:VAL:HG23	1.94	0.50
1:R:255:THR:HB	3:R:360:HOH:O	2.11	0.49
1:R:263:MET:HG3	1:R:292:PHE:CZ	2.47	0.49
1:O:183:ASP:HB2	1:O:192:ARG:NH2	2.27	0.49
1:Q:228:LEU:HD22	1:Q:245:LEU:HD11	1.95	0.49
1:O:80:ALA:HB1	1:O:109:HIS:CE1	2.47	0.49
1:Q:217:ILE:HD11	1:Q:221:ILE:HD12	1.95	0.49
1:P:183:ASP:HB2	1:P:192:ARG:NH2	2.27	0.48
1:O:83:LEU:HD13	1:O:85:TRP:CZ2	2.47	0.48
1:P:300:MET:CE	1:P:302:VAL:HG23	2.43	0.48
1:P:1:MET:C	1:P:1:MET:SD	2.92	0.48
1:O:263:MET:HG3	1:O:292:PHE:CZ	2.48	0.48
1:R:83:LEU:HD13	1:R:85:TRP:CZ2	2.48	0.48
1:P:251:LYS:HG2	1:P:254:VAL:HG23	1.96	0.48
1:P:263:MET:HG3	1:P:292:PHE:CZ	2.49	0.47
1:O:183:ASP:OD1	1:O:198:ARG:NH1	2.44	0.47
1:Q:242:LEU:HD11	1:Q:311:ALA:HB1	1.96	0.47
1:P:226:GLY:O	1:O:300:MET:HE3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:138:GLU:HG3	3:O:364:HOH:O	2.12	0.47
1:O:18:PHE:O	1:O:22:GLN:HG2	2.15	0.47
1:Q:245:LEU:HG	1:Q:247:VAL:HG13	1.96	0.47
1:R:57:GLU:H	1:R:68:ASN:ND2	2.13	0.46
1:O:242:LEU:HD11	1:O:311:ALA:HB1	1.97	0.46
1:P:20:ARG:O	1:P:24:VAL:HG23	2.16	0.46
1:O:80:ALA:HA	1:O:83:LEU:HD12	1.97	0.46
1:P:263:MET:HG3	1:P:292:PHE:CE1	2.50	0.46
1:Q:63:GLY:O	1:Q:73:LYS:HA	2.16	0.46
1:Q:263:MET:HG3	1:Q:292:PHE:CE1	2.51	0.46
1:Q:216:ALA:O	1:Q:220:VAL:HG23	2.16	0.46
1:R:242:LEU:HD11	1:R:311:ALA:HB1	1.98	0.46
1:O:34:ASP:OD2	1:O:36:THR:HG22	2.15	0.46
1:P:34:ASP:OD2	1:P:36:THR:HG22	2.16	0.45
1:R:224:ILE:HA	1:R:227:LYS:HD2	1.99	0.45
1:R:245:LEU:HG	1:R:247:VAL:HG13	1.98	0.45
1:O:46:LYS:HG3	3:O:349:HOH:O	2.17	0.45
1:P:16:LEU:CD1	1:P:317:GLU:HB3	2.47	0.45
1:O:245:LEU:HG	1:O:247:VAL:HG13	1.99	0.44
1:Q:300:MET:HE2	1:Q:302:VAL:HG23	1.99	0.44
1:O:105:LYS:HB2	3:O:360:HOH:O	2.18	0.44
1:P:243:THR:O	1:P:311:ALA:HA	2.18	0.44
1:Q:18:PHE:O	1:Q:22:GLN:HG2	2.18	0.44
1:Q:201:ALA:CB	3:Q:366:HOH:O	2.65	0.43
1:P:205:ASN:ND2	1:O:281:SER:CB	2.81	0.43
1:O:243:THR:O	1:O:311:ALA:HA	2.18	0.43
1:O:16:LEU:CD1	1:O:317:GLU:HB3	2.47	0.43
1:O:283:VAL:HG11	1:O:313:TRP:HB3	2.01	0.43
1:P:167:PHE:HD1	1:P:251:LYS:HD3	1.83	0.43
1:O:16:LEU:HD23	1:O:19:ARG:HD2	1.99	0.43
1:Q:47:TYR:HB3	1:O:282:ASP:OD1	2.19	0.43
1:P:109:HIS:HB2	1:P:117:VAL:HG21	1.99	0.43
1:O:79:ASP:CG	1:O:80:ALA:H	2.21	0.43
1:P:20:ARG:NH1	1:P:23:GLU:OE1	2.51	0.43
1:Q:205:ASN:HD22	1:R:281:SER:H	1.65	0.43
1:R:167:PHE:HD1	1:R:251:LYS:HD3	1.84	0.42
1:R:186:THR:HG22	1:O:180:TYR:CE1	2.54	0.42
1:Q:91:ASP:HA	1:Q:115:LYS:HD2	2.02	0.42
1:O:167:PHE:HD1	1:O:251:LYS:HD3	1.85	0.42
1:P:315:ASP:O	1:P:319:SER:HB2	2.19	0.42
1:P:31:ALA:HA	1:P:73:LYS:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:53:ARG:CG	1:Q:53:ARG:HH11	2.20	0.42
1:Q:309:LYS:HE3	3:Q:358:HOH:O	2.20	0.42
1:R:187:GLN:O	1:R:188:ASP:C	2.59	0.41
1:O:189:ALA:O	1:O:199:ALA:HB1	2.20	0.41
1:R:183:ASP:OD1	1:R:198:ARG:NH1	2.52	0.41
1:Q:281:SER:CB	1:R:205:ASN:HD21	2.33	0.41
1:P:245:LEU:HG	1:P:247:VAL:HG13	2.01	0.41
1:O:283:VAL:O	1:O:315:ASP:HB2	2.20	0.41
1:P:228:LEU:HD22	1:P:245:LEU:HD11	2.02	0.41
1:O:280:SER:O	1:O:283:VAL:HG22	2.21	0.41
1:P:37:ASP:OD2	1:P:39:ASP:HB3	2.21	0.41
1:R:334:LEU:HB3	3:R:367:HOH:O	2.21	0.41
1:R:178:HIS:HB3	1:R:234:ARG:HD3	2.03	0.41
1:P:309:LYS:HB2	1:O:173:LEU:HD13	2.03	0.41
1:Q:62:ASP:C	1:Q:64:GLY:H	2.24	0.41
1:Q:235:VAL:HA	1:Q:236:PRO:HD3	1.89	0.41
1:Q:263:MET:HG3	1:Q:292:PHE:CZ	2.56	0.41
1:O:31:ALA:HA	1:O:73:LYS:O	2.21	0.41
1:O:224:ILE:HA	1:O:227:LYS:HD2	2.02	0.41
1:P:174:MET:HE1	1:P:211:THR:HG21	2.02	0.41
1:P:189:ALA:O	1:P:199:ALA:HB1	2.20	0.40
1:R:198:ARG:HH21	1:R:209:ASN:HD22	1.69	0.40
1:Q:27:LEU:HD21	1:Q:329:ALA:HA	2.03	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:79:ASP:OD1	1:R:265:ASN:OD1[2_555]	2.02	0.18
1:Q:1:MET:N	1:Q:301:SER:OG[2_655]	2.06	0.14
1:Q:22:GLN:O	1:P:89:ASN:OD1[1_655]	2.10	0.10
1:P:111:GLU:OE1	1:P:261:GLU:OE1[2_556]	2.10	0.10
1:P:215:LYS:NZ	1:O:108:ALA:O[2_656]	2.10	0.10
1:Q:82:LYS:NZ	1:R:261:GLU:OE2[2_555]	2.11	0.09
1:R:110:ILE:CG2	1:R:257:GLU:CG[2_545]	2.17	0.03
1:Q:124:THR:OG1	1:R:269:GLU:OE2[2_555]	2.18	0.02

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	O	331/336 (98%)	312 (94%)	18 (5%)	1 (0%)	46	50
1	P	332/336 (99%)	319 (96%)	13 (4%)	0	100	100
1	Q	332/336 (99%)	314 (95%)	16 (5%)	2 (1%)	30	29
1	R	332/336 (99%)	319 (96%)	12 (4%)	1 (0%)	46	50
All	All	1327/1344 (99%)	1264 (95%)	59 (4%)	4 (0%)	46	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	R	80	ALA
1	O	68	ASN
1	Q	201	ALA
1	Q	61	VAL

### 5.3.2 Protein sidechains ⓘ

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	O	271/274 (99%)	261 (96%)	10 (4%)	41	50
1	P	272/274 (99%)	256 (94%)	16 (6%)	24	27
1	Q	272/274 (99%)	263 (97%)	9 (3%)	45	56
1	R	272/274 (99%)	261 (96%)	11 (4%)	38	47
All	All	1087/1096 (99%)	1041 (96%)	46 (4%)	36	44

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	Q	20	ARG
1	Q	53	ARG
1	Q	82	LYS
1	Q	85	TRP
1	Q	86	LYS
1	Q	211	THR
1	Q	239	THR
1	Q	269	GLU
1	Q	286	MET
1	P	1	MET
1	P	10	PHE
1	P	20	ARG
1	P	53	ARG
1	P	55	THR
1	P	62	ASP
1	P	82	LYS
1	P	85	TRP
1	P	86	LYS
1	P	115	LYS
1	P	166	ASP
1	P	215	LYS
1	P	239	THR
1	P	269	GLU
1	P	286	MET
1	P	334	LEU
1	R	55	THR
1	R	76	SER
1	R	82	LYS
1	R	85	TRP
1	R	86	LYS
1	R	115	LYS
1	R	166	ASP
1	R	269	GLU
1	R	286	MET
1	R	314	TYR
1	R	334	LEU
1	O	1	MET
1	O	62	ASP
1	O	68	ASN
1	O	82	LYS
1	O	85	TRP
1	O	86	LYS

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Mol	Chain	Res	Type
1	O	166	ASP
1	O	239	THR
1	O	269	GLU
1	O	286	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	Q	68	ASN
1	Q	133	ASN
1	Q	205	ASN
1	Q	209	ASN
1	Q	265	ASN
1	Q	296	GLN
1	Q	316	ASN
1	P	68	ASN
1	P	133	ASN
1	P	205	ASN
1	P	209	ASN
1	P	265	ASN
1	P	296	GLN
1	P	316	ASN
1	R	68	ASN
1	R	133	ASN
1	R	205	ASN
1	R	209	ASN
1	R	265	ASN
1	R	296	GLN
1	R	316	ASN
1	O	22	GLN
1	O	133	ASN
1	O	205	ASN
1	O	209	ASN
1	O	265	ASN
1	O	296	GLN
1	O	316	ASN

### 5.3.3 RNA ⓘ

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](#)

4 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	O	337	-	4,4,4	0.46	0	6,6,6	0.28	0
2	PO4	P	337	-	4,4,4	0.43	0	6,6,6	0.27	0
2	PO4	Q	337	-	4,4,4	0.50	0	6,6,6	0.31	0
2	PO4	R	337	-	4,4,4	0.61	0	6,6,6	0.27	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	O	337	-	-	0/0/0/0	0/0/0/0
2	PO4	P	337	-	-	0/0/0/0	0/0/0/0
2	PO4	Q	337	-	-	0/0/0/0	0/0/0/0
2	PO4	R	337	-	-	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.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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	O	333/336 (99%)	0.67	43 (12%) 5 4	47, 64, 91, 106	0
1	P	334/336 (99%)	0.51	26 (7%) 16 15	47, 61, 81, 97	0
1	Q	334/336 (99%)	0.61	33 (9%) 9 8	46, 61, 82, 92	0
1	R	334/336 (99%)	0.47	23 (6%) 20 19	46, 60, 82, 99	0
All	All	1335/1344 (99%)	0.57	125 (9%) 11 10	46, 61, 86, 106	0

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	Q	1	MET	7.0
1	O	112	ALA	7.0
1	R	304	ASP	6.1
1	R	55	THR	5.8
1	O	303	GLY	5.6
1	P	253	ASP	5.5
1	R	1	MET	5.4
1	R	124	THR	5.0
1	O	81	SER	5.0
1	O	252	GLN	4.9
1	Q	215	LYS	4.8
1	O	1	MET	4.8
1	P	304	ASP	4.7
1	P	60	VAL	4.6
1	R	334	LEU	4.5
1	Q	60	VAL	4.4
1	Q	192	ARG	4.3
1	Q	69	GLY	4.3
1	P	192	ARG	4.3
1	Q	59	GLU	4.3
1	Q	81	SER	4.2

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Mol	Chain	Res	Type	RSRZ
1	P	61	VAL	4.2
1	R	99	PHE	4.1
1	Q	63	GLY	4.1
1	Q	2	ALA	4.1
1	P	1	MET	4.1
1	P	215	LYS	4.0
1	P	66	ARG	3.9
1	Q	64	GLY	3.9
1	P	62	ASP	3.8
1	O	333	GLU	3.8
1	O	141	GLY	3.7
1	O	253	ASP	3.6
1	P	216	ALA	3.6
1	R	257	GLU	3.5
1	O	82	LYS	3.5
1	O	108	ALA	3.5
1	P	57	GLU	3.4
1	Q	66	ARG	3.4
1	Q	58	VAL	3.4
1	Q	107	GLN	3.4
1	O	60	VAL	3.3
1	O	106	ALA	3.3
1	O	89	ASN	3.2
1	O	37	ASP	3.1
1	Q	62	ASP	3.1
1	O	113	GLY	3.0
1	R	303	GLY	2.9
1	O	3	VAL	2.9
1	O	99	PHE	2.8
1	R	81	SER	2.8
1	Q	193	LYS	2.8
1	O	86	LYS	2.8
1	P	261	GLU	2.8
1	O	57	GLU	2.8
1	O	126	ASP	2.7
1	O	251	LYS	2.7
1	Q	242	LEU	2.7
1	O	62	ASP	2.7
1	Q	65	PHE	2.7
1	O	2	ALA	2.7
1	O	80	ALA	2.6
1	P	305	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	P	2	ALA	2.6
1	O	90	ILE	2.6
1	Q	86	LYS	2.6
1	O	104	ASP	2.6
1	P	258	GLN	2.6
1	O	35	LEU	2.6
1	R	56	GLY	2.5
1	O	304	ASP	2.5
1	Q	138	GLU	2.5
1	P	89	ASN	2.5
1	O	114	ALA	2.5
1	Q	133	ASN	2.5
1	O	254	VAL	2.4
1	R	2	ALA	2.4
1	R	126	ASP	2.4
1	R	192	ARG	2.4
1	Q	175	THR	2.4
1	P	213	ALA	2.4
1	O	124	THR	2.4
1	P	35	LEU	2.4
1	R	100	TYR	2.4
1	O	66	ARG	2.4
1	Q	82	LYS	2.4
1	O	245	LEU	2.4
1	O	192	ARG	2.4
1	Q	231	GLY	2.4
1	P	39	ASP	2.4
1	O	36	THR	2.4
1	R	252	GLN	2.4
1	O	111	GLU	2.3
1	R	251	LYS	2.3
1	R	331	LEU	2.3
1	O	258	GLN	2.3
1	P	99	PHE	2.3
1	R	104	ASP	2.3
1	P	142	SER	2.3
1	R	105	LYS	2.3
1	Q	216	ALA	2.2
1	P	228	LEU	2.2
1	P	254	VAL	2.2
1	O	250	GLU	2.2
1	Q	211	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	O	31	ALA	2.2
1	O	302	VAL	2.1
1	Q	127	LEU	2.1
1	Q	144	THR	2.1
1	Q	111	GLU	2.1
1	R	35	LEU	2.1
1	Q	87	ASP	2.1
1	O	71	GLU	2.1
1	Q	57	GLU	2.1
1	R	305	ARG	2.1
1	Q	212	GLY	2.1
1	P	249	LEU	2.1
1	P	140	ASP	2.1
1	Q	261	GLU	2.0
1	P	230	GLY	2.0
1	O	110	ILE	2.0
1	Q	251	LYS	2.0
1	R	82	LYS	2.0
1	R	39	ASP	2.0
1	O	307	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	PO4	O	337	5/5	0.97	0.13	-0.42	67,71,72,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	PO4	Q	337	5/5	0.96	0.12	-0.94	64,70,76,80	0
2	PO4	P	337	5/5	0.97	0.11	-1.88	53,68,71,72	0
2	PO4	R	337	5/5	0.98	0.08	-2.04	51,53,58,63	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.