



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2023 – 08:34 AM EDT

PDB ID : 1KC7
Title : Pyruvate Phosphate Dikinase with Bound Mg-phosphonopyruvate
Authors : Chen, C.C.; Howard, A.; Herzberg, O.
Deposited on : 2001-11-07
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35
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

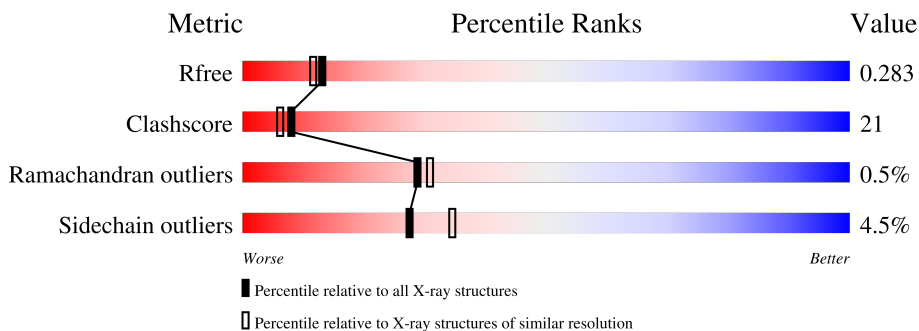
1 Overall quality at a glance

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}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (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 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	873	

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
4	PPR	A	1000	-	X	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6783 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 pyruvate phosphate dikinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	872	6752	4250	1143	1308	51	41	0	0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



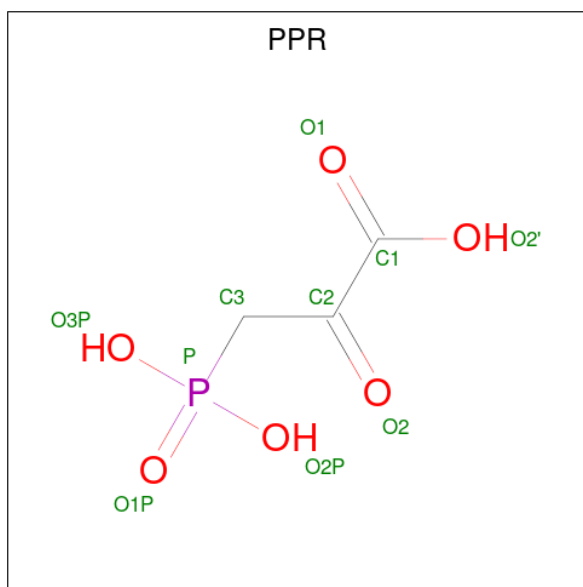
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 4 is PHOSPHONOPYRUVATE (three-letter code: PPR) (formula: $C_3H_5O_6P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O P 10 3 6 1	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	89.21Å 58.42Å 102.55Å 90.00° 94.99° 90.00°	Depositor
Resolution (Å)	25.00 – 2.20 36.17 – 1.73	Depositor EDS
% Data completeness (in resolution range)	96.3 (25.00-2.20) 90.0 (36.17-1.73)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.72 (at 1.73Å)	Xtrriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.198 , 0.269 0.240 , 0.283	Depositor DCC
R_{free} test set	4605 reflections (4.41%)	wwPDB-VP
Wilson B-factor (Å ²)	16.6	Xtrriage
Anisotropy	0.495	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6783	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: MG, PPR, 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	0.79	5/6876 (0.1%)	0.91	9/9269 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	812	VAL	CB-CG2	7.08	1.67	1.52
1	A	616	VAL	CB-CG1	6.42	1.66	1.52
1	A	749	ALA	CA-CB	-5.55	1.40	1.52
1	A	544	GLU	CG-CD	5.18	1.59	1.51
1	A	667	CYS	CB-SG	-5.12	1.73	1.81

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	537	ARG	NE-CZ-NH1	10.35	125.47	120.30
1	A	537	ARG	NE-CZ-NH2	-9.71	115.44	120.30
1	A	519	GLY	N-CA-C	-6.94	95.75	113.10
1	A	617	ARG	NE-CZ-NH2	-6.36	117.12	120.30
1	A	529	ASP	CB-CG-OD1	5.44	123.19	118.30
1	A	743	MET	CG-SD-CE	-5.34	91.66	100.20
1	A	754	ASP	CB-CG-OD1	5.21	122.99	118.30
1	A	145	MET	N-CA-C	-5.04	97.38	111.00
1	A	127	ARG	NE-CZ-NH2	-5.03	117.78	120.30

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	6752	0	6668	285	0
2	A	1	0	0	0	0
3	A	20	0	0	0	0
4	A	10	0	2	1	0
All	All	6783	0	6670	285	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 21.

All (285) 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:A:396:LEU:HB3	1:A:452:MET:HE3	1.28	1.15
1:A:172:THR:HG22	1:A:174:ASP:H	1.07	1.14
1:A:579:ILE:HD11	1:A:623:LEU:HD22	1.26	1.13
1:A:312:LEU:HD12	1:A:322:MET:HE2	1.32	1.08
1:A:396:LEU:HB3	1:A:452:MET:CE	1.84	1.06
1:A:557:ILE:HD13	1:A:609:LEU:HD21	1.50	0.94
1:A:172:THR:HG22	1:A:174:ASP:N	1.83	0.92
1:A:106:ILE:HD11	1:A:143:VAL:HG11	1.51	0.90
1:A:312:LEU:HD12	1:A:322:MET:CE	2.05	0.87
1:A:518:SER:O	1:A:521:PHE:HB3	1.77	0.85
1:A:309:ALA:HA	1:A:322:MET:CE	2.07	0.84
1:A:576:ARG:HD3	1:A:626:PHE:O	1.81	0.81
1:A:309:ALA:HA	1:A:322:MET:HE3	1.62	0.81
1:A:579:ILE:CD1	1:A:623:LEU:HD22	2.08	0.80
1:A:156:ILE:HG13	1:A:179:LEU:HD21	1.65	0.79
1:A:360:THR:OG1	1:A:363:GLU:HG3	1.84	0.77
1:A:165:VAL:HG13	1:A:170:ASP:HB2	1.68	0.76
1:A:468:VAL:HG11	1:A:505:ILE:HD11	1.68	0.75
1:A:529:ASP:CG	1:A:863:ARG:HH21	1.90	0.75
1:A:133:TYR:O	1:A:137:ILE:HG12	1.86	0.75
1:A:678:LYS:HG2	1:A:720:VAL:CG1	2.17	0.75
1:A:73:GLU:HG2	1:A:79:LYS:HA	1.69	0.74
1:A:223:TYR:HD2	1:A:436:ILE:HD11	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:556:GLY:C	1:A:557:ILE:HD12	2.09	0.73
1:A:743:MET:C	1:A:744:ILE:HD12	2.09	0.72
1:A:440:HIS:HB2	1:A:463:MET:HE2	1.71	0.71
1:A:317:ARG:HH22	1:A:363:GLU:CD	1.94	0.70
1:A:167:PHE:HB3	1:A:169:THR:HG23	1.72	0.70
1:A:189:GLU:HG3	1:A:189:GLU:O	1.90	0.70
1:A:396:LEU:HD12	1:A:451:GLY:HA2	1.72	0.70
1:A:505:ILE:N	1:A:505:ILE:HD12	2.06	0.70
1:A:4:TRP:H	1:A:64:GLN:HE22	1.40	0.70
1:A:580:LEU:HD23	1:A:651:LYS:HG2	1.75	0.68
1:A:425:ILE:HG21	1:A:496:ILE:HD11	1.75	0.68
1:A:381:PHE:CE1	1:A:510:ILE:HB	2.29	0.68
1:A:223:TYR:HD2	1:A:436:ILE:CD1	2.06	0.68
1:A:529:ASP:OD2	1:A:863:ARG:NH2	2.28	0.67
1:A:223:TYR:CD2	1:A:436:ILE:HD11	2.29	0.67
1:A:678:LYS:HB3	1:A:724:VAL:HG21	1.77	0.67
1:A:705:ILE:HD12	1:A:705:ILE:N	2.11	0.66
1:A:32:LEU:HD22	1:A:311:LYS:NZ	2.09	0.66
1:A:268:ILE:HD11	1:A:305:PHE:HE2	1.60	0.66
1:A:754:ASP:O	1:A:822:THR:HG21	1.96	0.66
1:A:628:PRO:HG2	1:A:634:GLN:HG2	1.76	0.65
1:A:151:HIS:O	1:A:155:ILE:HG12	1.97	0.65
1:A:4:TRP:H	1:A:64:GLN:NE2	1.94	0.65
1:A:557:ILE:HD13	1:A:609:LEU:CD2	2.25	0.64
1:A:165:VAL:CG1	1:A:170:ASP:HB2	2.28	0.64
1:A:599:LYS:HE2	1:A:686:GLU:HB3	1.79	0.64
1:A:172:THR:HB	1:A:175:ASP:OD2	1.98	0.64
1:A:644:THR:HG23	1:A:647:GLU:OE1	1.97	0.64
1:A:309:ALA:HA	1:A:322:MET:HE1	1.78	0.64
1:A:80:PHE:HA	1:A:87:LEU:HB3	1.79	0.64
1:A:96:ARG:HD2	1:A:233:TRP:CZ3	2.34	0.63
1:A:482:THR:HG22	1:A:491:ALA:HA	1.80	0.63
1:A:627:VAL:CG2	1:A:652:VAL:HG22	2.28	0.63
1:A:317:ARG:NH2	1:A:363:GLU:OE2	2.31	0.63
1:A:221:ILE:HD13	1:A:224:ARG:NH2	2.13	0.63
1:A:519:GLY:O	1:A:520:SER:HB2	1.99	0.63
1:A:343:ALA:HB3	1:A:344:PRO:HD3	1.81	0.63
1:A:782:ALA:HB1	1:A:786:LEU:HD22	1.81	0.62
1:A:62:GLN:HE21	1:A:205:MET:HG3	1.64	0.62
1:A:434:GLU:OE1	1:A:434:GLU:N	2.21	0.62
1:A:65:ILE:HG21	1:A:205:MET:HE1	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:ARG:HD2	1:A:233:TRP:CE3	2.34	0.61
1:A:268:ILE:HD11	1:A:305:PHE:CE2	2.35	0.61
1:A:354:VAL:HG11	1:A:527:TRP:CH2	2.35	0.61
1:A:830:ILE:HG22	1:A:852:VAL:HA	1.81	0.61
1:A:627:VAL:HG21	1:A:652:VAL:HG22	1.83	0.60
1:A:436:ILE:HG23	1:A:437:GLU:N	2.15	0.60
1:A:520:SER:HA	1:A:523:ARG:HB2	1.83	0.59
1:A:744:ILE:HD12	1:A:744:ILE:N	2.17	0.59
1:A:268:ILE:HD12	1:A:306:MET:SD	2.42	0.59
1:A:707:LEU:HD23	1:A:746:ILE:HD11	1.85	0.59
1:A:495:TYR:CD2	1:A:509:ASP:HB2	2.37	0.58
1:A:162:GLU:C	1:A:164:GLY:H	2.05	0.58
1:A:689:ILE:O	1:A:693:GLU:HG3	2.02	0.58
1:A:447:THR:OG1	1:A:469:SER:HB2	2.04	0.58
1:A:817:LYS:O	1:A:821:GLN:HG3	2.03	0.58
1:A:131:ASP:O	1:A:134:ARG:HG3	2.03	0.58
1:A:156:ILE:HG13	1:A:179:LEU:CD2	2.34	0.58
1:A:209:LYS:O	1:A:213:ARG:HB2	2.03	0.58
1:A:406:LYS:HG2	1:A:495:TYR:CE1	2.38	0.58
1:A:65:ILE:HG21	1:A:205:MET:CE	2.34	0.58
1:A:38:GLN:O	1:A:241:THR:HG22	2.04	0.57
1:A:488:HIS:CG	1:A:489:THR:H	2.22	0.57
1:A:742:THR:HG23	1:A:744:ILE:HD11	1.87	0.57
1:A:632:GLU:O	1:A:635:ALA:HB3	2.05	0.57
1:A:644:THR:OG1	1:A:647:GLU:HG3	2.05	0.57
1:A:79:LYS:CB	1:A:82:ASP:HB2	2.35	0.56
1:A:505:ILE:N	1:A:505:ILE:CD1	2.69	0.56
1:A:293:GLN:O	1:A:296:ASN:HB2	2.05	0.56
1:A:354:VAL:HG21	1:A:527:TRP:CZ3	2.41	0.56
1:A:390:GLU:O	1:A:392:ILE:HG23	2.06	0.56
1:A:454:SER:O	1:A:458:VAL:HG13	2.05	0.56
1:A:514:GLU:HG2	1:A:515:ALA:O	2.05	0.56
1:A:130:TYR:CZ	1:A:177:LYS:HG3	2.41	0.56
1:A:259:ARG:NH1	1:A:266:LYS:HG3	2.21	0.56
1:A:185:ALA:O	1:A:189:GLU:HB3	2.06	0.56
1:A:520:SER:O	1:A:524:ILE:HG12	2.06	0.56
1:A:630:THR:HG22	1:A:632:GLU:H	1.71	0.56
1:A:379:PRO:HB2	1:A:512:THR:OG1	2.05	0.55
1:A:78:LYS:NZ	1:A:85:ASP:OD1	2.39	0.55
1:A:128:PHE:CZ	1:A:247:LYS:HG3	2.42	0.55
1:A:584:VAL:O	1:A:588:GLU:HG3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:617:ARG:NH2	4:A:1000:PPR:O3P	2.40	0.55
1:A:112:ASN:O	1:A:116:VAL:HG12	2.06	0.55
1:A:704:MET:SD	1:A:743:MET:HG2	2.47	0.55
1:A:325:THR:OG1	1:A:334:LEU:HD11	2.07	0.54
1:A:144:VAL:HG12	1:A:145:MET:CE	2.37	0.54
1:A:580:LEU:HD12	1:A:641:MET:SD	2.48	0.54
1:A:707:LEU:HA	1:A:746:ILE:CD1	2.38	0.54
1:A:565:MET:HB3	1:A:598:GLN:HG2	1.90	0.54
1:A:628:PRO:HB3	1:A:633:GLU:HG2	1.90	0.53
1:A:34:MET:CE	1:A:308:LEU:HD22	2.39	0.53
1:A:262:SER:HB3	1:A:400:PRO:HD3	1.88	0.53
1:A:188:LYS:C	1:A:190:ALA:H	2.11	0.53
1:A:704:MET:SD	1:A:764:SER:HB3	2.49	0.53
1:A:519:GLY:C	1:A:521:PHE:H	2.12	0.53
1:A:572:ILE:O	1:A:576:ARG:HG3	2.09	0.53
1:A:268:ILE:CD1	1:A:305:PHE:HE2	2.22	0.52
1:A:109:LEU:HA	1:A:136:PHE:CE1	2.44	0.52
1:A:34:MET:HE2	1:A:308:LEU:HD22	1.92	0.52
1:A:678:LYS:HG2	1:A:720:VAL:HG13	1.90	0.52
1:A:689:ILE:HD12	1:A:734:SER:OG	2.10	0.52
1:A:113:ASP:OD1	1:A:184:LYS:NZ	2.42	0.52
1:A:681:THR:OG1	1:A:721:VAL:HG13	2.10	0.52
1:A:414:ALA:HB1	1:A:424:VAL:HG11	1.91	0.51
1:A:34:MET:SD	1:A:312:LEU:HD21	2.51	0.51
1:A:579:ILE:HD12	1:A:623:LEU:HD13	1.92	0.51
1:A:573:MET:CE	1:A:576:ARG:HH21	2.24	0.51
1:A:79:LYS:HB3	1:A:82:ASP:HB2	1.93	0.51
1:A:274:ILE:HD11	1:A:297:ASP:OD2	2.11	0.51
1:A:635:ALA:HA	1:A:645:LEU:HD13	1.93	0.51
1:A:322:MET:CE	1:A:333:PHE:HE2	2.24	0.50
1:A:757:ALA:HB3	1:A:822:THR:OG1	2.10	0.50
1:A:312:LEU:CD1	1:A:322:MET:HE2	2.23	0.50
1:A:7:LYS:HB2	1:A:10:GLU:HG3	1.93	0.50
1:A:189:GLU:O	1:A:189:GLU:CG	2.58	0.50
1:A:317:ARG:O	1:A:367:ARG:NH1	2.41	0.50
1:A:491:ALA:N	1:A:494:ASP:OD2	2.37	0.50
1:A:392:ILE:HG13	1:A:485:LEU:HD23	1.94	0.50
1:A:856:PRO:O	1:A:859:VAL:HG23	2.12	0.50
1:A:222:VAL:HG12	1:A:223:TYR:N	2.26	0.50
1:A:373:LEU:HD11	1:A:864:LEU:HD22	1.93	0.50
1:A:573:MET:HE2	1:A:576:ARG:HE	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:708:VAL:O	1:A:749:ALA:HB2	2.13	0.49
1:A:133:TYR:CE1	1:A:137:ILE:HD11	2.48	0.49
1:A:423:ARG:HH22	1:A:509:ASP:CG	2.15	0.49
1:A:742:THR:HG23	1:A:744:ILE:CD1	2.42	0.49
1:A:787:ASP:OD2	1:A:791:LYS:HE3	2.12	0.49
1:A:381:PHE:HE1	1:A:510:ILE:HB	1.77	0.49
1:A:830:ILE:HD13	1:A:831:CYS:N	2.28	0.49
1:A:744:ILE:HD11	1:A:763:PHE:CD1	2.48	0.49
1:A:672:THR:O	1:A:674:PRO:HD3	2.13	0.49
1:A:32:LEU:HD22	1:A:311:LYS:HZ2	1.76	0.48
1:A:711:LYS:HD3	1:A:755:ALA:O	2.14	0.48
1:A:763:PHE:CD2	1:A:826:LEU:HD21	2.48	0.48
1:A:276:ALA:HB2	1:A:286:ARG:HH22	1.78	0.48
1:A:504:LYS:C	1:A:505:ILE:HD12	2.33	0.48
1:A:722:VAL:HG23	1:A:740:ILE:HD12	1.95	0.48
1:A:714:LEU:HD21	1:A:760:ALA:HB2	1.96	0.48
1:A:223:TYR:HB2	1:A:436:ILE:CD1	2.43	0.48
1:A:375:GLN:HA	1:A:378:HIS:CE1	2.48	0.48
1:A:258:THR:O	1:A:259:ARG:HG2	2.13	0.48
1:A:415:LYS:O	1:A:419:GLU:HG2	2.13	0.48
1:A:223:TYR:CD2	1:A:436:ILE:CD1	2.92	0.48
1:A:573:MET:HE1	1:A:576:ARG:HH21	1.79	0.47
1:A:289:GLN:HE21	1:A:293:GLN:CD	2.17	0.47
1:A:548:ASN:O	1:A:551:LYS:HB3	2.13	0.47
1:A:575:ILE:O	1:A:579:ILE:HG12	2.15	0.47
1:A:782:ALA:HA	1:A:785:PHE:CE2	2.49	0.47
1:A:551:LYS:HB2	1:A:551:LYS:HE3	1.64	0.47
1:A:689:ILE:HG21	1:A:732:LYS:HD3	1.96	0.47
1:A:482:THR:HA	1:A:490:PHE:O	2.15	0.47
1:A:3:LYS:HD2	1:A:3:LYS:HA	1.65	0.46
1:A:667:CYS:SG	1:A:717:VAL:HG21	2.55	0.46
1:A:295:GLU:HB2	1:A:302:TYR:CD2	2.50	0.46
1:A:137:ILE:HD12	1:A:184:LYS:HG3	1.97	0.46
1:A:744:ILE:HD13	1:A:763:PHE:HB3	1.97	0.46
1:A:37:PRO:HB3	1:A:241:THR:HG23	1.97	0.46
1:A:162:GLU:C	1:A:164:GLY:N	2.67	0.46
1:A:396:LEU:HD12	1:A:451:GLY:CA	2.44	0.46
1:A:519:GLY:O	1:A:520:SER:CB	2.63	0.46
1:A:707:LEU:HA	1:A:746:ILE:HD11	1.98	0.46
1:A:373:LEU:N	1:A:373:LEU:HD23	2.30	0.46
1:A:130:TYR:CD1	1:A:177:LYS:HA	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:537:ARG:HD3	1:A:851:TYR:CD1	2.51	0.46
1:A:722:VAL:HG23	1:A:740:ILE:CD1	2.45	0.46
1:A:521:PHE:CZ	1:A:525:MET:HG3	2.51	0.46
1:A:79:LYS:HB2	1:A:82:ASP:HB2	1.98	0.45
1:A:82:ASP:O	1:A:86:PRO:HB3	2.15	0.45
1:A:572:ILE:O	1:A:576:ARG:CG	2.64	0.45
1:A:546:THR:HG23	1:A:557:ILE:HD11	1.99	0.45
1:A:122:LYS:O	1:A:122:LYS:HD3	2.17	0.45
1:A:148:PRO:O	1:A:150:SER:N	2.49	0.45
1:A:188:LYS:HA	1:A:194:GLU:O	2.16	0.45
1:A:552:LEU:HA	1:A:552:LEU:HD23	1.70	0.45
1:A:830:ILE:HB	1:A:849:LEU:CD2	2.47	0.45
1:A:102:MET:HG3	1:A:223:TYR:CD2	2.52	0.45
1:A:599:LYS:O	1:A:603:LYS:HG3	2.17	0.45
1:A:786:LEU:O	1:A:789:TYR:HB2	2.16	0.45
1:A:82:ASP:OD2	1:A:83:THR:N	2.50	0.45
1:A:468:VAL:HG22	1:A:503:GLY:HA2	1.99	0.45
1:A:271:GLU:HA	1:A:289:GLN:O	2.17	0.45
1:A:557:ILE:HD12	1:A:557:ILE:N	2.32	0.45
1:A:381:PHE:HZ	1:A:497:SER:HB3	1.81	0.44
1:A:797:SER:CB	1:A:802:ARG:HD3	2.47	0.44
1:A:244:PHE:HB3	1:A:246:ASN:OD1	2.18	0.44
1:A:436:ILE:HG23	1:A:437:GLU:H	1.80	0.44
1:A:483:PHE:CE1	1:A:490:PHE:HB2	2.53	0.44
1:A:165:VAL:HG12	1:A:166:HIS:N	2.33	0.44
1:A:618:TYR:CE1	1:A:703:ILE:HG23	2.53	0.44
1:A:144:VAL:HG12	1:A:145:MET:HE2	2.00	0.44
1:A:424:VAL:CG1	1:A:425:ILE:N	2.80	0.44
1:A:518:SER:C	1:A:519:GLY:O	2.50	0.44
1:A:677:ALA:O	1:A:678:LYS:C	2.55	0.44
1:A:688:ALA:O	1:A:692:LYS:HB2	2.17	0.44
1:A:574:LYS:HD2	1:A:574:LYS:N	2.33	0.43
1:A:61:ILE:O	1:A:65:ILE:HG12	2.18	0.43
1:A:80:PHE:CE2	1:A:200:PRO:HB2	2.53	0.43
1:A:29:MET:HA	1:A:32:LEU:HD12	1.99	0.43
1:A:68:ALA:O	1:A:71:TRP:HB3	2.19	0.43
1:A:436:ILE:CG2	1:A:437:GLU:N	2.81	0.43
1:A:230:PRO:HD2	1:A:233:TRP:CE2	2.53	0.43
1:A:518:SER:O	1:A:521:PHE:CB	2.58	0.43
1:A:57:ILE:HD11	1:A:209:LYS:HG2	2.01	0.43
1:A:707:LEU:CD2	1:A:746:ILE:HD11	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:830:ILE:HB	1:A:849:LEU:HD21	2.01	0.43
1:A:121:LYS:HA	1:A:121:LYS:HD3	1.56	0.43
1:A:59:GLN:CD	1:A:59:GLN:H	2.22	0.43
1:A:398:ALA:HB1	1:A:457:ALA:CB	2.49	0.43
1:A:499:ASP:OD2	1:A:501:SER:HB2	2.19	0.43
1:A:579:ILE:CD1	1:A:623:LEU:HD13	2.49	0.43
1:A:192:ASN:HD22	1:A:192:ASN:HA	1.66	0.43
1:A:757:ALA:HB3	1:A:822:THR:HG1	1.84	0.43
1:A:808:VAL:O	1:A:808:VAL:HG12	2.19	0.43
1:A:42:VAL:HG21	1:A:65:ILE:HD12	2.01	0.42
1:A:689:ILE:HD13	1:A:736:MET:HE3	2.01	0.42
1:A:188:LYS:HG3	1:A:194:GLU:O	2.20	0.42
1:A:685:MET:O	1:A:689:ILE:HG12	2.19	0.42
1:A:85:ASP:OD1	1:A:122:LYS:HG3	2.19	0.42
1:A:623:LEU:HD23	1:A:623:LEU:HA	1.82	0.42
1:A:633:GLU:HA	1:A:636:GLU:HG3	2.02	0.42
1:A:353:LEU:HD23	1:A:353:LEU:HA	1.82	0.42
1:A:399:SER:HA	1:A:400:PRO:HD2	1.84	0.42
1:A:717:VAL:O	1:A:721:VAL:HG23	2.20	0.42
1:A:173:ALA:O	1:A:177:LYS:HB2	2.20	0.42
1:A:460:ALA:HB1	1:A:465:THR:O	2.20	0.42
1:A:108:ASN:HB2	1:A:136:PHE:HB2	2.01	0.42
1:A:93:SER:HB3	1:A:211:VAL:HG13	2.02	0.41
1:A:99:MET:HG2	1:A:223:TYR:CE1	2.55	0.41
1:A:148:PRO:O	1:A:149:LYS:C	2.57	0.41
1:A:243:VAL:HG11	1:A:332:TYR:CG	2.55	0.41
1:A:7:LYS:HE2	1:A:71:TRP:CD1	2.55	0.41
1:A:284:GLY:HA2	1:A:288:PRO:HD3	2.02	0.41
1:A:765:PHE:CZ	1:A:816:VAL:HG23	2.55	0.41
1:A:514:GLU:HG2	1:A:515:ALA:N	2.35	0.41
1:A:563:GLU:CG	1:A:621:PRO:HD3	2.51	0.41
1:A:667:CYS:O	1:A:671:VAL:HG23	2.20	0.41
1:A:393:GLY:HA3	1:A:485:LEU:HD21	2.03	0.41
1:A:419:GLU:C	1:A:421:GLY:H	2.24	0.41
1:A:498:LEU:HD23	1:A:505:ILE:HG23	2.02	0.41
1:A:707:LEU:HA	1:A:746:ILE:HD13	2.01	0.41
1:A:165:VAL:CG1	1:A:166:HIS:N	2.83	0.41
1:A:342:THR:HB	1:A:344:PRO:HD2	2.02	0.41
1:A:148:PRO:C	1:A:150:SER:N	2.72	0.41
1:A:350:ALA:O	1:A:354:VAL:HG23	2.21	0.41
1:A:557:ILE:O	1:A:557:ILE:HG22	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:591:LEU:HD23	1:A:591:LEU:HA	1.84	0.41
1:A:763:PHE:CE2	1:A:826:LEU:HD21	2.55	0.41
1:A:427:VAL:HA	1:A:446:LEU:O	2.22	0.41
1:A:570:ASP:OD2	1:A:571:ARG:NE	2.53	0.41
1:A:674:PRO:O	1:A:678:LYS:HG3	2.21	0.41
1:A:399:SER:O	1:A:500:GLY:HA3	2.21	0.40
1:A:689:ILE:HD13	1:A:736:MET:SD	2.61	0.40
1:A:144:VAL:HG12	1:A:145:MET:HE1	2.03	0.40
1:A:521:PHE:CE2	1:A:525:MET:HG3	2.57	0.40
1:A:630:THR:HG22	1:A:632:GLU:N	2.34	0.40
1:A:3:LYS:HE2	1:A:5:VAL:O	2.22	0.40
1:A:277:GLN:O	1:A:278:GLY:C	2.60	0.40
1:A:460:ALA:O	1:A:461:ARG:C	2.58	0.40
1:A:429:LEU:HD11	1:A:449:ARG:HD3	2.03	0.40
1:A:519:GLY:C	1:A:521:PHE:N	2.75	0.40
1:A:859:VAL:N	1:A:860:PRO:CD	2.84	0.40

There are no symmetry-related clashes.

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	870/873 (100%)	820 (94%)	46 (5%)	4 (0%)	29 31

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	GLU
1	A	190	ALA
1	A	149	LYS
1	A	856	PRO

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	713/714 (100%)	681 (96%)	32 (4%)	27 34

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	SER
1	A	24	CYS
1	A	31	ILE
1	A	98	SER
1	A	121	LYS
1	A	127	ARG
1	A	146	GLU
1	A	156	ILE
1	A	157	ASP
1	A	169	THR
1	A	213	ARG
1	A	216	ASP
1	A	268	ILE
1	A	279	GLU
1	A	300	ASP
1	A	322	MET
1	A	458	VAL
1	A	473	GLU
1	A	488	HIS
1	A	505	ILE
1	A	513	GLN
1	A	517	VAL
1	A	520	SER
1	A	576	ARG
1	A	578	MET
1	A	636	GLU
1	A	698	ASP
1	A	743	MET
1	A	747	PRO
1	A	780	ASP

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Mol	Chain	Res	Type
1	A	786	LEU
1	A	830	ILE

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

Mol	Chain	Res	Type
1	A	56	GLN
1	A	62	GLN
1	A	64	GLN
1	A	138	GLN
1	A	192	ASN
1	A	217	ASN
1	A	440	HIS
1	A	513	GLN
1	A	624	HIS
1	A	727	GLN
1	A	845	HIS

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](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 for Mogul analysis.

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	SO4	A	1003	-	4,4,4	1.55	1 (25%)	6,6,6	1.57	1 (16%)
4	PPR	A	1000	2	9,9,9	5.49	8 (88%)	12,13,13	2.72	6 (50%)
3	SO4	A	1002	-	4,4,4	0.25	0	6,6,6	0.46	0
3	SO4	A	1004	-	4,4,4	0.36	0	6,6,6	0.12	0
3	SO4	A	1005	-	4,4,4	0.45	0	6,6,6	0.29	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
4	PPR	A	1000	2	-	3/8/9/9	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1000	PPR	P-C3	10.36	1.96	1.79
4	A	1000	PPR	C3-C2	-9.86	1.36	1.51
4	A	1000	PPR	O2-C2	4.39	1.32	1.23
4	A	1000	PPR	P-O2P	-4.00	1.45	1.54
4	A	1000	PPR	P-O3P	-3.05	1.47	1.54
4	A	1000	PPR	P-O1P	2.83	1.56	1.50
4	A	1000	PPR	O1-C1	2.66	1.29	1.22
4	A	1000	PPR	O2'-C1	-2.65	1.22	1.30
3	A	1003	SO4	O2-S	-2.41	1.32	1.46

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1000	PPR	O1-C1-C2	-4.74	115.40	121.72
4	A	1000	PPR	C3-C2-C1	4.45	130.30	118.14
4	A	1000	PPR	O2-C2-C1	-3.98	113.67	119.43
3	A	1003	SO4	O4-S-O3	3.24	122.88	109.06
4	A	1000	PPR	O2'-C1-C2	3.20	122.72	113.97
4	A	1000	PPR	O3P-P-O2P	2.54	115.49	108.08
4	A	1000	PPR	O2-C2-C3	-2.41	116.02	120.27

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1000	PPR	C2-C3-P-O2P
4	A	1000	PPR	C2-C3-P-O3P
4	A	1000	PPR	C2-C3-P-O1P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1000	PPR	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](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.