



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:49 PM UTC

PDB ID : 1QE5 / pdb_00001qe5
Title : PURINE NUCLEOSIDE PHOSPHORYLASE FROM CELLULOMONAS SP.
IN COMPLEX WITH PHOSPHATE
Authors : Tebbe, J.; Bzowska, A.; Wielgus-Kutrowska, B.; Schroeder, W.; Kazimierczuk,
Z.; Shugar, D.; Saenger, W.; Koellner, G.
Deposited on : 1999-07-13
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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

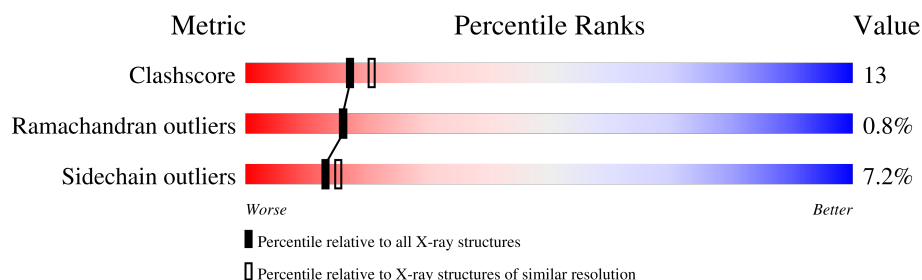
1 Overall quality at a glance

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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	266	 71% 27% .
1	B	266	 72% 24% .
1	C	266	 77% 18% . .

2 Entry composition [i](#)

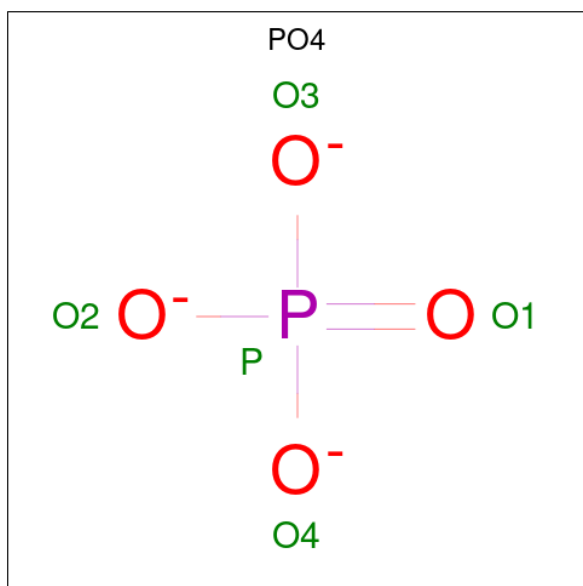
There are 4 unique types of molecules in this entry. The entry contains 6092 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PENTOSYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	0	0
			1938	1213	348	372	5			
1	B	266	Total	C	N	O	S	0	0	0
			1938	1213	348	372	5			
1	C	266	Total	C	N	O	S	0	0	0
			1938	1213	348	372	5			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Ca 1	0	0

- Molecule 4 is water.

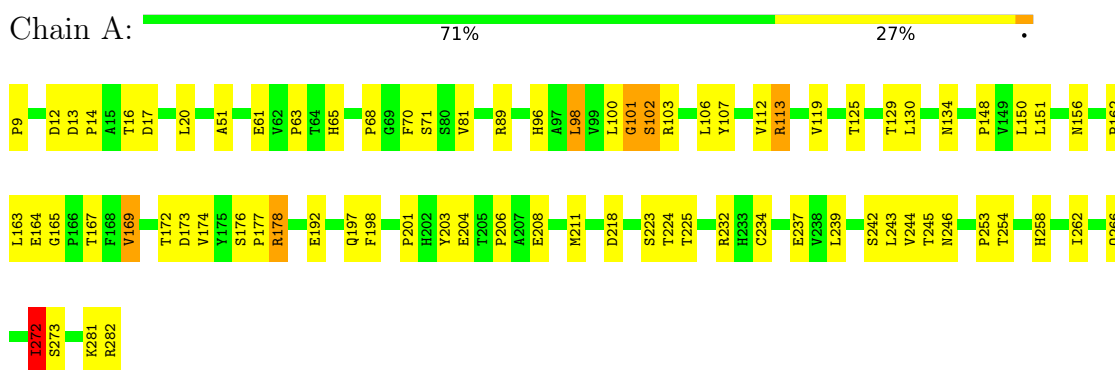
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	81	Total 81	O 81	0	0
4	B	95	Total 95	O 95	0	0
4	C	86	Total 86	O 86	0	0

3 Residue-property plots

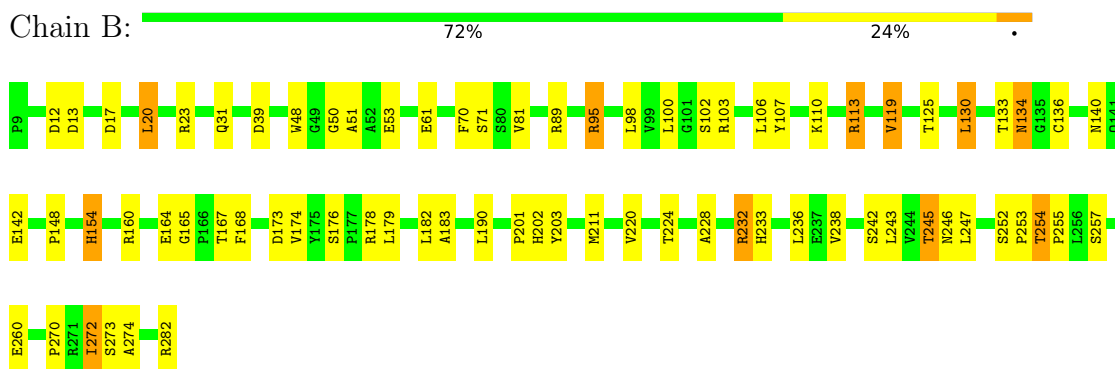
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

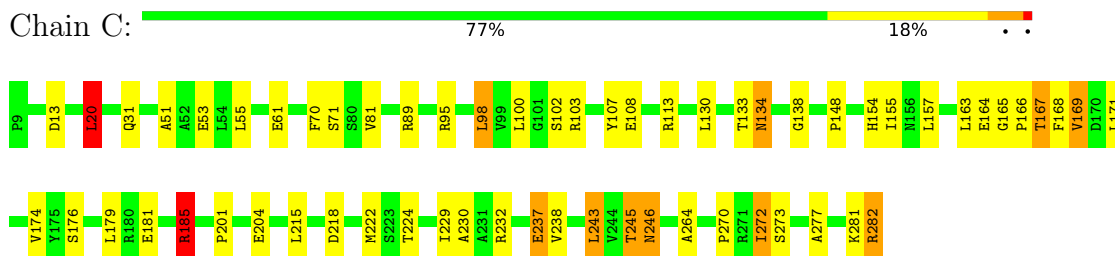
• Molecule 1: PENTOSYLTRANSFERASE



• Molecule 1: PENTOSYLTRANSFERASE



• Molecule 1: PENTOSYLTRANSFERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.12Å 108.90Å 119.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.00 – 2.20	Depositor
% Data completeness (in resolution range)	86.3 (28.00-2.20)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.200 , 0.260	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6092	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, CA

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	1.15	4/1977 (0.2%)	1.37	13/2700 (0.5%)
1	B	1.07	1/1977 (0.1%)	1.31	10/2700 (0.4%)
1	C	1.13	3/1977 (0.2%)	1.33	15/2700 (0.6%)
All	All	1.12	8/5931 (0.1%)	1.34	38/8100 (0.5%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	272	ILE	CA-CB	7.54	1.63	1.54
1	B	119	VAL	CA-C	6.60	1.60	1.52
1	C	243	LEU	CA-C	5.79	1.59	1.52
1	C	133	THR	CA-C	5.45	1.59	1.52
1	A	239	LEU	C-O	-5.31	1.18	1.24
1	C	230	ALA	CA-C	5.30	1.59	1.52
1	A	224	THR	CA-CB	5.24	1.62	1.53
1	A	206	PRO	CA-C	5.07	1.60	1.52

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	125	THR	N-CA-C	-8.64	102.74	113.28
1	A	254	THR	CA-C-N	8.43	128.93	119.83
1	A	254	THR	C-N-CA	8.43	128.93	119.83
1	A	224	THR	N-CA-C	7.79	127.40	110.80
1	B	165	GLY	CA-C-N	7.71	129.48	119.84
1	B	165	GLY	C-N-CA	7.71	129.48	119.84
1	B	238	VAL	N-CA-C	7.64	118.87	108.17
1	C	185	ARG	N-CA-C	-7.35	103.35	111.36
1	B	176	SER	N-CA-C	7.30	120.08	109.71
1	C	169	VAL	N-CA-C	6.95	118.07	108.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	PRO	N-CA-C	-6.72	104.18	113.53
1	B	245	THR	N-CA-C	-6.64	104.12	111.36
1	C	222	MET	N-CA-C	6.50	121.04	111.87
1	A	218	ASP	N-CA-C	-6.33	106.15	114.31
1	A	239	LEU	CA-C-N	-6.08	116.90	122.80
1	A	239	LEU	C-N-CA	-6.08	116.90	122.80
1	B	224	THR	N-CA-C	5.84	123.23	110.80
1	A	20	LEU	N-CA-C	-5.78	104.95	112.23
1	C	245	THR	N-CA-C	-5.70	105.07	111.28
1	B	183	ALA	N-CA-C	-5.68	104.99	111.07
1	A	198	PHE	CA-C-N	-5.64	113.08	119.28
1	A	198	PHE	C-N-CA	-5.64	113.08	119.28
1	C	108	GLU	N-CA-C	-5.61	106.44	113.28
1	C	218	ASP	N-CA-C	-5.52	106.55	113.72
1	B	274	ALA	N-CA-C	-5.50	105.36	111.36
1	C	264	ALA	N-CA-C	-5.42	105.45	111.36
1	C	167	THR	N-CA-C	5.40	117.47	108.02
1	C	238	VAL	N-CA-C	5.38	116.22	108.48
1	C	53	GLU	N-CA-C	5.35	118.03	111.82
1	C	165	GLY	CA-C-N	5.31	126.48	119.84
1	C	165	GLY	C-N-CA	5.31	126.48	119.84
1	B	190	LEU	N-CA-C	5.22	115.99	109.57
1	C	20	LEU	N-CA-C	-5.20	105.69	111.36
1	A	125	THR	N-CA-C	-5.17	106.95	113.20
1	C	224	THR	N-CA-C	5.13	121.72	110.80
1	A	244	VAL	CA-C-N	5.07	127.50	120.29
1	A	244	VAL	C-N-CA	5.07	127.50	120.29
1	C	237	GLU	N-CA-C	-5.01	101.54	109.96

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	A	1938	0	1931	56	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1938	0	1931	55	0
1	C	1938	0	1931	39	0
2	A	5	0	0	1	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
3	B	1	0	0	0	0
4	A	81	0	0	11	0
4	B	95	0	0	8	0
4	C	86	0	0	6	0
All	All	6092	0	5793	146	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (146) 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:B:270:PRO:HG2	4:B:565:HOH:O	1.54	1.05
1:A:113:ARG:HD2	1:A:164:GLU:OE2	1.63	0.99
1:B:103:ARG:HD2	4:B:531:HOH:O	1.67	0.94
1:B:103:ARG:CD	4:B:531:HOH:O	2.16	0.92
1:C:163:LEU:HD13	1:C:167:THR:HG21	1.55	0.87
1:B:254:THR:HG22	1:B:254:THR:O	1.74	0.86
1:A:129:THR:HG21	1:A:282:ARG:HH22	1.38	0.86
1:B:51:ALA:HA	1:B:273:SER:HB3	1.57	0.83
1:B:282:ARG:HH11	1:B:282:ARG:HG3	1.43	0.83
1:B:257:SER:OG	1:B:260:GLU:HG3	1.81	0.81
1:C:204:GLU:OE1	4:C:451:HOH:O	1.97	0.81
1:C:113:ARG:HD2	1:C:164:GLU:OE2	1.81	0.80
1:C:232:ARG:HD3	4:C:409:HOH:O	1.85	0.76
1:A:232:ARG:HD3	4:A:544:HOH:O	1.87	0.75
1:C:181:GLU:OE1	1:C:185:ARG:NH2	2.20	0.72
1:A:165:GLY:O	1:A:167:THR:N	2.23	0.70
1:B:110:LYS:HB3	4:B:420:HOH:O	1.92	0.70
1:B:70:PHE:O	1:B:71:SER:HB3	1.91	0.70
1:A:262:ILE:O	1:A:266:GLN:HG3	1.92	0.69
1:A:129:THR:HG21	1:A:282:ARG:NH2	2.08	0.68
1:B:113:ARG:HD2	1:B:164:GLU:OE2	1.94	0.67
1:B:254:THR:O	1:B:254:THR:CG2	2.41	0.67
1:C:245:THR:O	1:C:246:ASN:HB3	1.93	0.67
1:B:140:ASN:OD1	1:B:142:GLU:HG3	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:GLU:HB3	1:A:81:VAL:HG21	1.77	0.66
1:B:107:TYR:HB2	1:B:201:PRO:HD3	1.78	0.65
1:C:89:ARG:HD2	1:C:95:ARG:HG2	1.76	0.65
1:B:103:ARG:HD3	4:B:531:HOH:O	1.89	0.65
1:B:61:GLU:HB3	1:B:81:VAL:HG21	1.77	0.65
1:B:148:PRO:HD2	4:B:502:HOH:O	1.98	0.63
1:A:163:LEU:HD11	1:A:169:VAL:HG11	1.80	0.62
1:B:282:ARG:HH11	1:B:282:ARG:CG	2.09	0.62
1:A:282:ARG:CD	4:A:323:HOH:O	2.47	0.62
1:A:282:ARG:NE	4:A:323:HOH:O	2.32	0.61
1:A:61:GLU:HB3	1:A:81:VAL:CG2	2.31	0.61
1:A:96:HIS:HD2	4:A:448:HOH:O	1.84	0.60
1:B:61:GLU:HB3	1:B:81:VAL:CG2	2.31	0.60
1:C:281:LYS:O	1:C:282:ARG:HB2	2.01	0.59
1:A:96:HIS:CD2	4:A:448:HOH:O	2.56	0.58
1:B:211:MET:HG3	1:C:215:LEU:HD22	1.84	0.58
1:A:89:ARG:NH1	4:A:519:HOH:O	2.36	0.58
1:A:63:PRO:HB2	1:A:65:HIS:CE1	2.39	0.58
1:A:262:ILE:CG2	1:A:266:GLN:HE21	2.15	0.58
1:C:107:TYR:HB2	1:C:201:PRO:HD3	1.85	0.58
1:C:148:PRO:HB3	1:C:272:ILE:HD11	1.86	0.58
1:C:51:ALA:HA	1:C:273:SER:HB3	1.86	0.57
1:B:134:ASN:HD22	1:B:134:ASN:C	2.13	0.57
1:A:258:HIS:NE2	1:A:262:ILE:HD11	2.20	0.56
1:C:61:GLU:HB3	1:C:81:VAL:CG2	2.36	0.56
1:A:98:LEU:HD13	1:A:100:LEU:HD11	1.88	0.55
1:B:245:THR:O	1:B:246:ASN:HB3	2.06	0.55
1:C:61:GLU:HB3	1:C:81:VAL:HG21	1.87	0.55
1:C:176:SER:HB2	1:C:237:GLU:HG3	1.89	0.54
1:A:148:PRO:HB3	1:A:272:ILE:HD11	1.88	0.54
1:A:70:PHE:O	1:A:71:SER:HB3	2.07	0.54
1:B:89:ARG:HD2	1:B:95:ARG:HG2	1.89	0.54
1:B:174:VAL:HA	1:B:232:ARG:HD3	1.90	0.54
1:B:203:TYR:O	1:C:171:LEU:HD12	2.08	0.54
1:C:89:ARG:HD3	1:C:95:ARG:NH1	2.24	0.53
1:B:253:PRO:C	1:B:255:PRO:HD3	2.34	0.53
1:C:174:VAL:HA	1:C:232:ARG:HD2	1.89	0.53
1:A:17:ASP:C	1:A:17:ASP:OD1	2.49	0.53
1:B:282:ARG:CG	1:B:282:ARG:NH1	2.67	0.52
1:C:89:ARG:HD2	1:C:95:ARG:CG	2.38	0.52
1:B:253:PRO:O	1:B:255:PRO:HD3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:THR:O	1:A:246:ASN:HB3	2.09	0.51
1:C:154:HIS:CD2	1:C:154:HIS:C	2.88	0.51
1:B:273:SER:HB3	4:B:490:HOH:O	2.10	0.51
1:B:50:GLY:O	1:B:53:GLU:HG2	2.11	0.50
1:B:107:TYR:CD1	1:B:107:TYR:C	2.89	0.50
1:A:281:LYS:HE2	4:A:489:HOH:O	2.10	0.50
1:A:281:LYS:O	1:A:282:ARG:HB2	2.12	0.50
1:B:228:ALA:O	1:B:232:ARG:HG2	2.12	0.50
1:A:150:LEU:O	1:A:192:GLU:HA	2.12	0.49
1:B:134:ASN:C	1:B:134:ASN:ND2	2.70	0.49
1:C:134:ASN:HD22	1:C:134:ASN:C	2.19	0.49
1:B:39:ASP:HB3	1:B:95:ARG:NH1	2.27	0.49
1:B:100:LEU:HD12	1:B:100:LEU:N	2.28	0.49
1:C:98:LEU:HD22	1:C:100:LEU:HD12	1.94	0.49
1:A:112:VAL:HG21	1:A:162:PRO:HA	1.94	0.48
1:C:167:THR:CG2	1:C:169:VAL:HG13	2.44	0.48
1:C:13:ASP:OD1	1:C:13:ASP:C	2.56	0.48
1:A:178:ARG:NH2	1:A:237:GLU:OE1	2.31	0.48
1:B:247:LEU:HD12	1:B:252:SER:HB2	1.95	0.47
1:A:156:ASN:HB3	4:C:547:HOH:O	2.14	0.47
1:C:20:LEU:C	1:C:20:LEU:HD13	2.40	0.47
1:C:134:ASN:C	1:C:134:ASN:ND2	2.73	0.47
1:A:174:VAL:HA	1:A:232:ARG:HD2	1.97	0.47
1:C:107:TYR:C	1:C:107:TYR:CD1	2.93	0.47
1:A:163:LEU:CD1	1:A:169:VAL:HG11	2.44	0.47
1:A:107:TYR:HB2	1:A:201:PRO:HD3	1.97	0.46
1:A:197:GLN:NE2	1:A:225:THR:OG1	2.49	0.46
1:A:282:ARG:HD3	4:A:323:HOH:O	2.11	0.46
1:C:179:LEU:HD11	1:C:237:GLU:HG2	1.98	0.46
1:C:185:ARG:HH11	1:C:185:ARG:HA	1.81	0.46
1:B:89:ARG:HD2	1:B:95:ARG:CG	2.45	0.46
1:B:232:ARG:NE	1:B:232:ARG:HA	2.30	0.46
1:B:201:PRO:HG2	1:C:163:LEU:HB2	1.98	0.45
1:A:262:ILE:HG22	1:A:266:GLN:HE21	1.80	0.45
1:C:229:ILE:HD11	4:C:447:HOH:O	2.15	0.45
1:A:102:SER:OG	1:A:103:ARG:N	2.49	0.45
1:C:179:LEU:CD1	1:C:237:GLU:HG2	2.47	0.45
1:B:20:LEU:C	1:B:20:LEU:HD13	2.42	0.45
1:A:234:CYS:SG	4:A:436:HOH:O	2.37	0.44
1:B:160:ARG:NH2	4:B:406:HOH:O	2.49	0.44
1:B:154:HIS:CD2	1:B:154:HIS:C	2.95	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:PHE:O	1:C:71:SER:CB	2.65	0.44
1:C:113:ARG:CD	1:C:164:GLU:OE2	2.60	0.44
1:B:70:PHE:O	1:B:71:SER:CB	2.63	0.44
1:B:272:ILE:HA	1:B:272:ILE:HD12	1.78	0.44
1:B:182:LEU:HD13	1:B:282:ARG:HD2	1.99	0.44
1:A:148:PRO:HG2	4:A:510:HOH:O	2.18	0.44
1:A:101:GLY:O	1:A:102:SER:O	2.36	0.44
1:A:13:ASP:HA	1:A:14:PRO:HD3	1.91	0.43
1:B:48:TRP:CE2	1:B:133:THR:HB	2.53	0.43
1:A:119:VAL:HG23	4:A:326:HOH:O	2.16	0.43
1:C:138:GLY:HA3	4:C:391:HOH:O	2.19	0.43
1:A:100:LEU:O	1:A:101:GLY:O	2.37	0.42
1:B:179:LEU:HD23	1:B:179:LEU:HA	1.84	0.42
1:C:282:ARG:N	1:C:282:ARG:HD3	2.31	0.42
1:A:204:GLU:H	1:A:204:GLU:CD	2.26	0.42
1:B:17:ASP:OD1	1:B:17:ASP:C	2.63	0.42
1:A:148:PRO:HA	1:A:242:SER:O	2.19	0.42
1:A:204:GLU:CD	1:A:204:GLU:N	2.78	0.42
1:B:13:ASP:OD1	1:B:13:ASP:C	2.63	0.42
1:A:223:SER:OG	2:A:303:PO4:O1	2.35	0.42
1:A:98:LEU:HD13	1:A:100:LEU:CD1	2.49	0.42
1:A:203:TYR:CE2	1:B:168:PHE:HB3	2.55	0.42
1:B:136:CYS:HB2	1:B:220:VAL:O	2.19	0.42
1:A:282:ARG:N	1:A:282:ARG:HD2	2.32	0.42
1:B:202:HIS:HD2	1:B:203:TYR:O	2.03	0.42
1:C:270:PRO:HG2	4:C:464:HOH:O	2.20	0.42
1:C:55:LEU:HD12	1:C:277:ALA:HA	2.02	0.42
1:A:51:ALA:HA	1:A:273:SER:HB3	2.02	0.41
1:A:208:GLU:O	1:A:211:MET:HB3	2.20	0.41
1:C:155:ILE:HG22	1:C:157:LEU:HG	2.02	0.41
1:A:9:PRO:HB2	1:A:16:THR:HG23	2.01	0.41
1:A:172:THR:O	1:A:172:THR:HG22	2.20	0.41
1:A:165:GLY:C	1:A:167:THR:N	2.78	0.41
1:A:63:PRO:HB2	1:A:65:HIS:HE1	1.85	0.41
1:A:258:HIS:NE2	1:A:262:ILE:CD1	2.83	0.41
1:B:167:THR:HG21	1:B:233:HIS:CD2	2.56	0.41
1:A:176:SER:HA	1:A:177:PRO:HD3	1.94	0.40
1:B:20:LEU:HD22	1:B:20:LEU:O	2.21	0.40
1:B:81:VAL:O	1:B:81:VAL:HG13	2.22	0.40
1:B:119:VAL:HG22	1:B:130:LEU:HD11	2.04	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	262/266 (98%)	255 (97%)	5 (2%)	2 (1%)	16	16
1	B	262/266 (98%)	250 (95%)	12 (5%)	0	100	100
1	C	262/266 (98%)	252 (96%)	6 (2%)	4 (2%)	8	6
All	All	786/798 (98%)	757 (96%)	23 (3%)	6 (1%)	16	16

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	102	SER
1	A	101	GLY
1	C	102	SER
1	C	168	PHE
1	C	166	PRO
1	C	246	ASN

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	199/199 (100%)	186 (94%)	13 (6%)	15	18
1	B	199/199 (100%)	179 (90%)	20 (10%)	7	7
1	C	199/199 (100%)	189 (95%)	10 (5%)	22	28
All	All	597/597 (100%)	554 (93%)	43 (7%)	13	15

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ASP
1	A	68	PRO
1	A	98	LEU
1	A	106	LEU
1	A	113	ARG
1	A	130	LEU
1	A	134	ASN
1	A	151	LEU
1	A	169	VAL
1	A	173	ASP
1	A	178	ARG
1	A	243	LEU
1	A	272	ILE
1	B	12	ASP
1	B	20	LEU
1	B	23	ARG
1	B	31	GLN
1	B	95	ARG
1	B	98	LEU
1	B	102	SER
1	B	106	LEU
1	B	113	ARG
1	B	130	LEU
1	B	134	ASN
1	B	154	HIS
1	B	173	ASP
1	B	178	ARG
1	B	232	ARG
1	B	236	LEU
1	B	242	SER
1	B	243	LEU
1	B	254	THR
1	B	272	ILE
1	C	20	LEU
1	C	31	GLN
1	C	98	LEU
1	C	103	ARG
1	C	130	LEU
1	C	134	ASN
1	C	185	ARG
1	C	243	LEU
1	C	272	ILE

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Mol	Chain	Res	Type
1	C	282	ARG

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	28	HIS
1	A	31	GLN
1	A	197	GLN
1	A	266	GLN
1	B	65	HIS
1	B	134	ASN
1	B	197	GLN
1	C	65	HIS
1	C	134	ASN
1	C	197	GLN
1	C	266	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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
2	PO4	B	302	-	4,4,4	0.81	0	6,6,6	1.30	0
2	PO4	A	303	-	4,4,4	2.17	3 (75%)	6,6,6	0.99	0
2	PO4	C	301	-	4,4,4	0.84	0	6,6,6	1.28	0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	303	PO4	P-O1	2.95	1.57	1.50
2	A	303	PO4	P-O2	-2.32	1.47	1.54
2	A	303	PO4	P-O3	-2.03	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	303	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1
1	A	1
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	71:SER	C	80:SER	N	9.74
1	A	71:SER	C	80:SER	N	9.66
1	C	71:SER	C	80:SER	N	8.65

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.