



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 06:12 AM GMT

PDB ID : 4E6X
Title : ClbP in complex boron-based inhibitor
Authors : Cougnoux, A.; Delmas, J.; Bonnet, R.
Deposited on : 2012-03-16
Resolution : 2.24 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

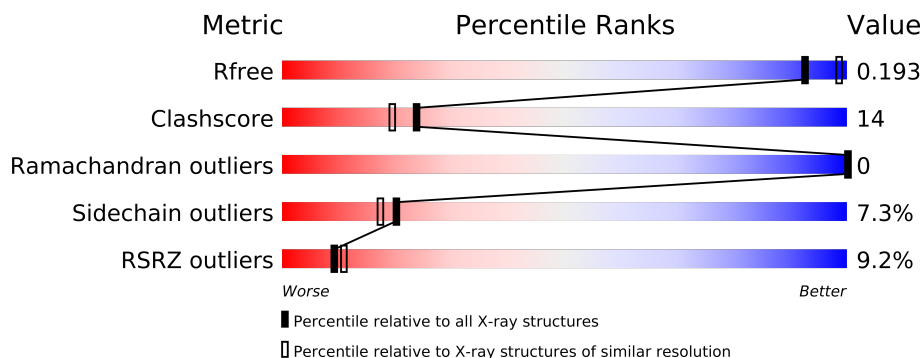
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.24 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}	66092	1112 (2.26-2.22)
Clashscore	79885	1317 (2.26-2.22)
Ramachandran outliers	78287	1282 (2.26-2.22)
Sidechain outliers	78261	1282 (2.26-2.22)
RSRZ outliers	66119	1112 (2.26-2.22)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	336	
1	B	336	
1	C	336	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	PO4	A	401	-	X
3	ONG	B	400	-	X

2 Entry composition i

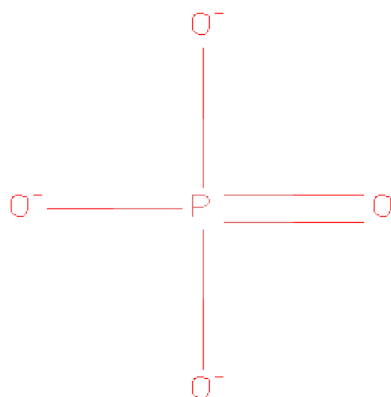
There are 4 unique types of molecules in this entry. The entry contains 7261 atoms, of which 0 are hydrogen and 0 are deuterium.

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 ClbP peptidase.

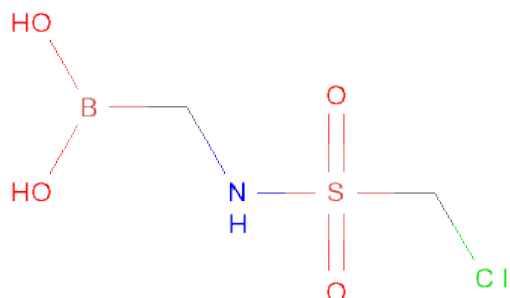
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	0	2	0
			2383	1516	415	443	9			
1	B	301	Total	C	N	O	S	0	0	0
			2321	1484	399	429	9			
1	C	305	Total	C	N	O	S	0	1	0
			2350	1501	405	435	9			

- Molecule 2 is PHOSPHATE ION (three-letter code: 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 ({[(CHLOROMETHYL)SULFONYL]AMINO}METHYL)BORONICACID (three-letter code: 0NG) (formula: C₂H₇BClNO₄S).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	B	1	Total	B	C	Cl	N	O	S	0	0
			10	1	2	1	1	4	1		
3	C	1	Total	B	C	Cl	N	O	S	0	0
			10	1	2	1	1	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	61	Total	O	0	0
			61	61		
4	B	49	Total	O	0	0
			49	49		
4	C	62	Total	O	0	0
			62	62		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	105.05Å 152.53Å 86.89Å 90.00° 123.43° 90.00°	Depositor
Resolution (Å)	52.55 – 2.24 52.55 – 2.24	Depositor EDS
% Data completeness (in resolution range)	79.0 (52.55-2.24) 78.0 (52.55-2.24)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 2.25Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.200 , 0.249 0.180 , 0.193	Depositor DCC
R_{free} test set	2195 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 26.7	EDS
Estimated twinning fraction	0.467 for 1/2*h+1/2*k+l,3/2*h-1/2*k+l,-l 0.459 for 1/2*h-1/2*k+l,-3/2*h-1/2*k-l,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 43180 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7261	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, 0NG

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	0.71	1/2435 (0.0%)	0.83	0/3308
1	B	0.60	0/2368	0.81	1/3217 (0.0%)
1	C	0.60	0/2400	0.79	0/3262
All	All	0.64	1/7203 (0.0%)	0.81	1/9787 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	337	CYS	CB-SG	-5.68	1.72	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	183	LYS	CB-CA-C	-5.13	100.13	110.40

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2383	0	2370	58	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2321	0	2321	75	0
1	C	2350	0	2341	69	0
2	A	5	0	0	1	0
2	B	5	0	0	1	0
2	C	5	0	0	1	0
3	B	10	0	5	0	0
3	C	10	0	5	0	0
4	A	61	0	0	0	0
4	B	49	0	0	0	0
4	C	62	0	0	0	0
All	All	7261	0	7042	202	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (202) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:152:PHE:HD1	1:C:152:PHE:C	1.52	1.12
1:A:283:LEU:HB3	1:A:284:PRO:HD2	1.29	1.08
1:C:134:LYS:HG2	1:C:135:PRO:HD2	1.30	1.07
1:B:344:GLN:O	1:B:344:GLN:HG3	1.57	1.04
1:C:152:PHE:CD1	1:C:152:PHE:C	2.29	1.02
1:C:245:LYS:O	1:C:245:LYS:HG2	1.55	1.02
1:C:152:PHE:O	1:C:152:PHE:HD1	1.49	0.94
1:A:243:PHE:N	1:A:243:PHE:HD1	1.67	0.92
1:C:134:LYS:HG2	1:C:135:PRO:CD	2.00	0.90
1:A:333:ASN:HD22	1:A:333:ASN:H	1.16	0.90
1:A:243:PHE:CD1	1:A:243:PHE:N	2.39	0.88
1:C:46:LEU:O	1:C:46:LEU:HD23	1.72	0.87
1:A:245:LYS:O	1:A:245:LYS:HD3	1.74	0.86
1:C:152:PHE:O	1:C:152:PHE:CD1	2.29	0.83
1:A:333:ASN:H	1:A:333:ASN:ND2	1.75	0.83
1:B:248:LEU:CD1	1:B:248:LEU:C	2.46	0.83
1:B:245:LYS:CD	1:B:245:LYS:O	2.29	0.80
1:A:172:ARG:HG2	1:A:173:ASN:N	1.96	0.79
1:C:46:LEU:CD2	1:C:46:LEU:O	2.30	0.78
1:B:245:LYS:CG	1:B:245:LYS:O	2.31	0.78
1:C:134:LYS:CG	1:C:135:PRO:HD2	2.10	0.78
1:A:283:LEU:HB3	1:A:284:PRO:CD	2.11	0.77
1:A:245:LYS:CG	1:A:245:LYS:O	2.32	0.77
1:B:90:VAL:HB	1:B:235:LYS:HA	1.66	0.76
1:C:90:VAL:HB	1:C:235:LYS:HA	1.67	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:258:VAL:HA	1:B:262:TYR:HB2	1.69	0.75
1:C:241:LEU:O	1:C:241:LEU:HD23	1.86	0.75
1:A:280:ARG:HH11	1:A:280:ARG:HG2	1.51	0.75
1:B:258:VAL:HG13	1:B:262:TYR:O	1.88	0.74
1:C:245:LYS:O	1:C:245:LYS:CG	2.34	0.74
1:A:309:ILE:HB	2:A:401:PO4:O1	1.87	0.73
1:B:146:HIS:CD2	1:B:187:ALA:H	2.07	0.73
1:B:220:MET:HE1	1:B:273:TRP:HE3	1.54	0.73
1:A:245:LYS:CD	1:A:245:LYS:O	2.36	0.72
1:B:146:HIS:HD2	1:B:187:ALA:H	1.37	0.72
1:B:89:THR:CG2	1:B:236:ALA:HB2	2.19	0.72
1:B:240:LYS:HZ2	1:B:248:LEU:HD23	1.54	0.71
1:C:46:LEU:CD2	1:C:46:LEU:C	2.59	0.71
1:A:146:HIS:CD2	1:A:187:ALA:H	2.08	0.71
1:B:240:LYS:NZ	1:B:248:LEU:HD23	2.07	0.70
1:B:245:LYS:HD3	1:B:245:LYS:O	1.92	0.70
1:C:280:ARG:O	1:C:288:ARG:HG3	1.92	0.69
1:C:240:LYS:O	1:C:246:PRO:HD2	1.92	0.69
1:A:97:SER:HB3	1:A:263:ILE:HD12	1.74	0.69
1:A:232:ILE:HG22	1:A:233:VAL:O	1.93	0.69
1:B:248:LEU:HD13	1:B:248:LEU:C	2.13	0.68
1:C:283:LEU:HB3	1:C:284:PRO:HD2	1.76	0.67
1:C:248:LEU:HB3	1:C:333:ASN:HD22	1.60	0.67
1:A:146:HIS:HD2	1:A:187:ALA:H	1.42	0.65
1:A:245:LYS:HG2	1:A:245:LYS:O	1.96	0.65
1:B:248:LEU:HD12	1:B:248:LEU:C	2.17	0.63
1:A:243:PHE:H	1:A:243:PHE:HD1	1.29	0.63
1:C:92:GLU:HG2	1:C:264:HIS:NE2	2.14	0.63
1:B:245:LYS:O	1:B:245:LYS:HG2	1.99	0.62
1:B:151:PRO:O	1:B:152:PHE:HB2	1.97	0.62
1:C:57:PRO:HD3	1:C:354:MET:HE3	1.81	0.62
1:A:119:ASP:O	1:A:122:THR:HG23	1.99	0.62
1:C:245:LYS:O	1:C:245:LYS:HE2	2.00	0.61
1:C:257:HIS:HD2	1:C:262:TYR:HE2	1.49	0.60
1:B:218:LEU:HD13	1:B:220:MET:CE	2.32	0.60
1:B:248:LEU:HD13	1:B:248:LEU:O	2.03	0.59
1:A:226:VAL:O	1:A:258:VAL:HG21	2.02	0.58
1:A:280:ARG:HG2	1:A:280:ARG:NH1	2.16	0.58
1:B:248:LEU:HD12	1:B:333:ASN:ND2	2.19	0.58
1:B:233:VAL:O	1:B:234:ASN:HB2	2.02	0.57
1:B:309:ILE:HB	2:B:401:PO4:O1	2.04	0.57
1:A:204:LYS:HE3	1:A:212:GLU:OE1	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:92:GLU:HG2	1:C:264:HIS:CD2	2.40	0.57
1:A:333:ASN:ND2	1:A:333:ASN:N	2.42	0.57
1:C:233:VAL:O	1:C:234:ASN:HB2	2.04	0.56
1:A:129:LEU:HD23	1:A:174:GLU:HB3	1.86	0.56
1:C:146:HIS:CD2	1:C:187:ALA:H	2.23	0.56
1:C:97:SER:HB3	1:C:263:ILE:HD12	1.88	0.56
1:B:46:LEU:HD23	1:B:46:LEU:O	2.05	0.56
1:B:143:PHE:HD1	1:B:190:ASN:HD22	1.55	0.55
1:A:293:ASN:HA	1:A:296:ARG:HE	1.71	0.55
1:A:226:VAL:HG23	1:A:226:VAL:O	2.07	0.55
1:B:218:LEU:HD13	1:B:220:MET:HE3	1.89	0.54
1:A:143:PHE:HD1	1:A:190:ASN:HD22	1.55	0.54
1:A:207:THR:HG23	1:A:225:ALA:O	2.07	0.54
1:C:46:LEU:HD23	1:C:46:LEU:C	2.26	0.54
1:B:220:MET:HE2	1:B:273:TRP:CA	2.38	0.54
1:A:333:ASN:HD22	1:A:333:ASN:N	1.85	0.54
1:C:248:LEU:HB3	1:C:333:ASN:ND2	2.22	0.54
1:C:368:ASN:HD22	1:C:371:ARG:HH21	1.56	0.54
1:B:238:GLY:HA3	1:B:333:ASN:HB2	1.89	0.54
1:A:64:THR:HB	1:A:69:ARG:HG3	1.90	0.53
1:C:257:HIS:HD2	1:C:262:TYR:CE2	2.26	0.53
1:C:240:LYS:HZ2	1:C:248:LEU:CD1	2.21	0.53
1:B:248:LEU:HD12	1:B:333:ASN:HD21	1.74	0.53
1:B:104:VAL:HG23	1:B:214:LEU:HD22	1.91	0.52
1:B:245:LYS:C	1:B:245:LYS:HD3	2.29	0.52
1:B:358:LEU:N	1:B:358:LEU:HD12	2.24	0.52
1:C:46:LEU:HD22	1:C:46:LEU:C	2.29	0.52
1:C:134:LYS:CG	1:C:135:PRO:CD	2.81	0.52
1:B:344:GLN:HG2	1:B:346:ILE:CD1	2.39	0.52
1:B:131:TYR:O	1:B:132:GLN:HB2	2.09	0.52
1:B:315:TRP:CE3	1:B:325:ILE:HG22	2.45	0.52
1:B:295:TRP:CZ3	1:B:325:ILE:HD13	2.45	0.51
1:C:56:VAL:HG12	1:C:354:MET:HB3	1.91	0.51
1:B:344:GLN:HG2	1:B:346:ILE:HD11	1.93	0.51
1:C:258:VAL:HA	1:C:262:TYR:HB2	1.92	0.51
1:A:258:VAL:HA	1:A:262:TYR:HB2	1.92	0.51
1:B:121:ILE:HG23	1:B:121:ILE:O	2.10	0.51
1:B:220:MET:HE2	1:B:273:TRP:HA	1.93	0.51
1:C:184:PHE:CE2	1:C:311:TYR:HB2	2.46	0.50
1:C:226:VAL:O	1:C:258:VAL:HG21	2.10	0.50
1:B:220:MET:HE1	1:B:273:TRP:CE3	2.41	0.50
1:B:283:LEU:HB3	1:B:284:PRO:HD2	1.93	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:89:THR:HG21	1:B:236:ALA:HB2	1.92	0.50
1:C:123:TYR:O	1:C:124:LEU:HD12	2.12	0.50
1:A:192:ASP:OD1	1:A:257:HIS:HA	2.12	0.50
1:A:218:LEU:HB2	1:A:220:MET:HE2	1.92	0.49
1:B:240:LYS:NZ	1:B:248:LEU:CD2	2.74	0.49
1:A:244:GLY:O	1:A:246:PRO:HD3	2.13	0.49
1:B:281:LYS:HA	1:B:288:ARG:HD2	1.95	0.49
1:A:78:ASP:OD2	1:A:81:SER:HB3	2.13	0.49
1:B:207:THR:HG23	1:B:225:ALA:O	2.13	0.49
1:C:209:VAL:O	1:C:213:ARG:HB2	2.13	0.49
1:C:309:ILE:HB	2:C:401:PO4:O2	2.13	0.49
1:C:46:LEU:O	1:C:46:LEU:HD22	2.12	0.48
1:C:134:LYS:HG2	1:C:135:PRO:N	2.28	0.48
1:B:335:SER:HB3	1:B:355:ASN:HA	1.95	0.48
1:B:220:MET:CE	1:B:273:TRP:HA	2.43	0.48
1:C:123:TYR:C	1:C:124:LEU:HD12	2.34	0.48
1:C:240:LYS:HZ2	1:C:248:LEU:CD2	2.27	0.48
1:A:56:VAL:HG12	1:A:354:MET:HB3	1.96	0.47
1:A:238:GLY:HA3	1:A:333:ASN:HB2	1.97	0.47
1:C:241:LEU:CD2	1:C:241:LEU:O	2.61	0.47
1:B:245:LYS:HD3	1:B:245:LYS:N	2.29	0.47
1:C:241:LEU:HG	1:C:242:GLY:O	2.15	0.47
1:C:266:THR:HB	1:C:268:PRO:HD2	1.96	0.47
1:B:238:GLY:HA3	1:B:333:ASN:CB	2.44	0.47
1:A:233:VAL:HG23	1:A:234:ASN:OD1	2.15	0.47
1:A:368:ASN:HD22	1:A:371:ARG:HH21	1.62	0.47
1:A:127:MET:O	1:A:127:MET:HG2	2.15	0.47
1:B:64:THR:O	1:B:346:ILE:HA	2.15	0.46
1:B:218:LEU:HD13	1:B:220:MET:HE1	1.97	0.46
1:C:333:ASN:N	1:C:333:ASN:OD1	2.48	0.46
1:B:247:VAL:CG2	1:B:248:LEU:N	2.78	0.46
1:B:333:ASN:N	1:B:333:ASN:OD1	2.48	0.46
1:A:95:SER:HG	1:A:98:LYS:HZ3	1.62	0.46
1:C:283:LEU:HB3	1:C:284:PRO:CD	2.44	0.46
1:C:368:ASN:ND2	1:C:371:ARG:HH21	2.14	0.46
1:C:267:LEU:HB3	1:C:268:PRO:HD3	1.97	0.46
1:B:204:LYS:HB3	1:B:208:GLU:OE1	2.16	0.46
1:B:103:LEU:O	1:B:107:ILE:HG13	2.16	0.46
1:A:95:SER:OG	1:A:98:LYS:NZ	2.44	0.46
1:B:232:ILE:HD12	1:B:232:ILE:N	2.31	0.45
1:B:103:LEU:HD23	1:B:214:LEU:HD11	1.99	0.45
1:B:245:LYS:H	1:B:245:LYS:HD3	1.81	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:172:ARG:HG2	1:A:173:ASN:H	1.76	0.45
1:C:64:THR:O	1:C:346:ILE:HA	2.16	0.45
1:B:361:GLN:HG3	1:B:361:GLN:O	2.16	0.45
1:A:191:TYR:O	1:A:259:PRO:HB2	2.17	0.45
1:A:104:VAL:HG23	1:A:214:LEU:HD11	1.98	0.45
1:A:64:THR:O	1:A:346:ILE:HA	2.16	0.44
1:A:95:SER:HB3	1:A:327:HIS:CE1	2.52	0.44
1:A:283:LEU:CB	1:A:284:PRO:HD2	2.18	0.44
1:B:238:GLY:C	1:B:333:ASN:HB3	2.38	0.44
1:B:130:ASN:OD1	1:B:135:PRO:HA	2.18	0.44
1:B:286:THR:HG23	1:B:287:LEU:N	2.33	0.44
1:B:132:GLN:HB3	1:B:133:GLY:H	1.47	0.44
1:C:213:ARG:HA	1:C:213:ARG:HD2	1.71	0.44
1:A:240:LYS:O	1:A:246:PRO:HD2	2.18	0.43
1:C:46:LEU:HD11	1:C:365:ASP:HB3	1.99	0.43
1:C:224:VAL:HG22	1:C:225:ALA:N	2.34	0.43
1:A:295:TRP:CE3	1:A:325:ILE:HD13	2.53	0.43
1:C:152:PHE:CD1	1:C:153:SER:N	2.85	0.43
1:B:211:ALA:HA	1:B:215:THR:OG1	2.18	0.43
1:A:48:HIS:O	1:A:52:GLN:HG3	2.17	0.43
1:B:214:LEU:HA	1:B:214:LEU:HD12	1.79	0.43
1:B:119:ASP:OD1	1:B:120:ILE:N	2.51	0.43
1:C:224:VAL:HG13	1:C:224:VAL:O	2.19	0.43
1:C:280:ARG:HB2	1:C:291:MET:SD	2.59	0.43
1:C:238:GLY:HA3	1:C:333:ASN:HB2	2.00	0.43
1:C:199:GLU:HG2	1:C:204:LYS:O	2.18	0.42
1:C:262:TYR:O	1:C:264:HIS:ND1	2.52	0.42
1:A:90:VAL:HB	1:A:235:LYS:HA	2.00	0.42
1:A:69:ARG:NH2	1:A:275:ASP:OD2	2.52	0.42
1:C:64:THR:HB	1:C:69:ARG:HG3	2.01	0.42
1:C:152:PHE:HD1	1:C:153:SER:N	2.11	0.42
1:B:195:GLY:HA3	1:B:259:PRO:HG2	2.01	0.42
1:A:233:VAL:O	1:A:234:ASN:HB2	2.18	0.42
1:A:108:LEU:HD13	1:A:201:VAL:HG11	2.02	0.42
1:B:90:VAL:N	1:B:234:ASN:O	2.25	0.42
1:B:134:LYS:H	1:B:134:LYS:HG3	1.68	0.42
1:C:57:PRO:HD2	1:C:353:ASN:OD1	2.20	0.42
1:B:46:LEU:HD23	1:B:46:LEU:C	2.39	0.42
1:A:213:ARG:N	1:A:213:ARG:HD2	2.34	0.41
1:B:111:GLU:CD	1:B:213:ARG:HH22	2.23	0.41
1:A:130:ASN:N	1:A:174:GLU:O	2.38	0.41
1:B:281:LYS:HA	1:B:288:ARG:CD	2.50	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:183:LYS:HB2	1:C:183:LYS:HE3	1.61	0.41
1:C:57:PRO:O	1:C:75:GLY:HA3	2.21	0.41
1:C:87:LEU:HD23	1:C:87:LEU:HA	1.93	0.41
1:A:245:LYS:HB2	1:A:354:MET:HE2	2.02	0.41
1:B:46:LEU:C	1:B:46:LEU:CD2	2.89	0.41
1:B:295:TRP:CE3	1:B:325:ILE:HD13	2.55	0.41
1:C:295:TRP:CZ3	1:C:325:ILE:HD13	2.56	0.40
1:C:100:PHE:HB2	1:C:263:ILE:CD1	2.51	0.40
1:B:134:LYS:HB3	1:B:135:PRO:HD2	2.02	0.40

There are no symmetry-related clashes.

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	A	304/336 (90%)	296 (97%)	8 (3%)	0	100	100
1	B	293/336 (87%)	281 (96%)	12 (4%)	0	100	100
1	C	298/336 (89%)	289 (97%)	9 (3%)	0	100	100
All	All	895/1008 (89%)	866 (97%)	29 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	A	251/276 (91%)	228 (91%)	23 (9%)	13	9
1	B	247/276 (90%)	232 (94%)	15 (6%)	26	25

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	249/276 (90%)	232 (93%)	17 (7%)	22	20
All	All	747/828 (90%)	692 (93%)	55 (7%)	20	17

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	81	SER
1	A	95	SER
1	A	115	ARG
1	A	122	THR
1	A	126	GLU
1	A	137	SER
1	A	172	ARG
1	A	176	LEU
1	A	213	ARG
1	A	227	LYS
1	A	241	LEU
1	A	243	PHE
1	A	245	LYS
1	A	247	VAL
1	A	256[A]	ASN
1	A	256[B]	ASN
1	A	257	HIS
1	A	296	ARG
1	A	333	ASN
1	A	344	GLN
1	A	358	LEU
1	A	370	LEU
1	B	41	GLU
1	B	49	GLN
1	B	121	ILE
1	B	126	GLU
1	B	128	ARG
1	B	176	LEU
1	B	214	LEU
1	B	233	VAL
1	B	243	PHE
1	B	245	LYS
1	B	247	VAL
1	B	248	LEU
1	B	333	ASN

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Mol	Chain	Res	Type
1	B	344	GLN
1	B	370	LEU
1	C	46	LEU
1	C	53	GLU
1	C	132	GLN
1	C	152	PHE
1	C	176	LEU
1	C	183	LYS
1	C	213	ARG
1	C	214	LEU
1	C	240	LYS
1	C	241	LEU
1	C	243	PHE
1	C	245	LYS
1	C	247	VAL
1	C	248	LEU
1	C	333	ASN
1	C	358	LEU
1	C	370	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	116	GLN
1	A	132	GLN
1	A	146	HIS
1	A	190	ASN
1	A	333	ASN
1	A	368	ASN
1	B	49	GLN
1	B	146	HIS
1	B	190	ASN
1	B	321	GLN
1	B	357	ASN
1	B	368	ASN
1	C	116	GLN
1	C	132	GLN
1	C	146	HIS
1	C	190	ASN
1	C	368	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	PO4	A	401	-	4,4,4	0.41	0	6,6,6	0.34	0
3	0NG	B	400	1	9,9,9	4.06	5 (55%)	11,12,12	3.63	5 (45%)
2	PO4	B	401	-	4,4,4	0.35	0	6,6,6	0.33	0
3	0NG	C	400	1	9,9,9	4.27	5 (55%)	11,12,12	3.70	6 (54%)
2	PO4	C	401	-	4,4,4	0.46	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
2	PO4	A	401	-	-	0/0/0/0	0/0/0/0
3	0NG	B	400	1	-	0/5/9/9	0/0/0/0
2	PO4	B	401	-	-	0/0/0/0	0/0/0/0
3	0NG	C	400	1	-	0/5/9/9	0/0/0/0
2	PO4	C	401	-	-	0/0/0/0	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	400	0NG	S08-N07	8.60	1.85	1.60
3	B	400	0NG	S08-N07	7.07	1.81	1.60
3	B	400	0NG	O09-S08	6.70	1.54	1.43
3	B	400	0NG	O10-S08	6.20	1.53	1.43
3	C	400	0NG	O09-S08	6.05	1.53	1.43
3	C	400	0NG	O10-S08	5.72	1.53	1.43
3	C	400	0NG	C11-S08	3.00	1.83	1.77
3	C	400	0NG	C06-N07	2.68	1.50	1.46
3	B	400	0NG	C06-N07	2.57	1.49	1.46
3	B	400	0NG	C11-S08	2.39	1.82	1.77

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	400	0NG	O04-B03-C06	-7.48	107.49	121.37
3	B	400	0NG	O04-B03-C06	-7.17	108.06	121.37
3	B	400	0NG	O05-B03-C06	-6.69	108.96	121.37
3	C	400	0NG	O05-B03-C06	-5.99	110.25	121.37
3	C	400	0NG	O09-S08-O10	-5.57	109.07	119.30
3	B	400	0NG	O09-S08-O10	-5.03	110.07	119.30
3	B	400	0NG	C06-N07-S08	-3.73	109.68	121.64
3	C	400	0NG	O10-S08-N07	3.41	115.35	107.07
3	B	400	0NG	O10-S08-N07	2.72	113.68	107.07
3	C	400	0NG	C06-N07-S08	-2.43	113.86	121.64
3	C	400	0NG	CL1-C11-S08	2.08	114.70	109.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 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	310/336 (92%)	0.81	28 (9%)	10 11	17, 35, 74, 97	0
1	B	301/336 (89%)	0.82	30 (9%)	8 9	19, 34, 74, 101	0
1	C	305/336 (90%)	0.77	27 (8%)	10 11	19, 33, 73, 97	0
All	All	916/1008 (90%)	0.80	85 (9%)	9 10	17, 34, 74, 101	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	244	GLY	16.0
1	A	249	PHE	11.4
1	C	242	GLY	10.8
1	A	244	GLY	10.5
1	A	243	PHE	10.2
1	A	173	ASN	9.9
1	C	243	PHE	9.8
1	A	242	GLY	9.7
1	C	153	SER	9.0
1	B	152	PHE	7.4
1	A	234	ASN	7.3
1	C	247	VAL	7.3
1	A	247	VAL	7.2
1	B	243	PHE	6.8
1	A	238	GLY	6.6
1	C	174	GLU	6.5
1	B	247	VAL	6.4
1	C	152	PHE	6.4
1	A	245	LYS	6.1
1	A	246	PRO	5.8
1	B	248	LEU	5.6
1	C	246	PRO	5.6
1	A	255	ARG	5.4

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Mol	Chain	Res	Type	RSRZ
1	B	246	PRO	5.2
1	B	245	LYS	4.9
1	A	41	GLU	4.7
1	A	257	HIS	4.6
1	B	151	PRO	4.5
1	B	262	TYR	4.2
1	A	133	GLY	4.0
1	C	267	LEU	4.0
1	C	241	LEU	4.0
1	B	261	ALA	3.9
1	C	259	PRO	3.9
1	B	242	GLY	3.9
1	C	245	LYS	3.9
1	C	244	GLY	3.8
1	B	237	SER	3.6
1	C	257	HIS	3.6
1	C	248	LEU	3.5
1	C	133	GLY	3.5
1	B	135	PRO	3.4
1	B	233	VAL	3.4
1	A	240	LYS	3.3
1	B	133	GLY	3.3
1	A	135	PRO	3.3
1	A	254	ALA	3.3
1	B	112	GLY	3.2
1	A	227	LYS	3.2
1	B	203	GLY	3.2
1	A	198	ILE	3.1
1	B	178	PHE	3.0
1	C	262	TYR	3.0
1	A	256[A]	ASN	3.0
1	C	256	ASN	3.0
1	C	203	GLY	2.9
1	A	129	LEU	2.8
1	A	231	ILE	2.8
1	C	225	ALA	2.8
1	B	129	LEU	2.8
1	B	232	ILE	2.7
1	C	177	LEU	2.6
1	B	130	ASN	2.6
1	C	240	LYS	2.6
1	A	172	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	332	PRO	2.4
1	A	171	LEU	2.4
1	A	226	VAL	2.4
1	A	248	LEU	2.3
1	C	151	PRO	2.3
1	C	120	ILE	2.2
1	C	132	GLN	2.2
1	B	277	TRP	2.2
1	B	241	LEU	2.2
1	B	121	ILE	2.2
1	B	80	ALA	2.1
1	B	258	VAL	2.1
1	A	121	ILE	2.1
1	B	132	GLN	2.1
1	A	132	GLN	2.0
1	B	176	LEU	2.0
1	C	140	VAL	2.0
1	C	232	ILE	2.0
1	C	126	GLU	2.0
1	B	119	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	PO4	A	401	5/5	0.27	4.79	26,27,31,35	5

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	0NG	B	400	10/10	0.28	2.12	37,43,49,57	10
3	0NG	C	400	10/10	0.22	1.69	41,50,57,66	10
2	PO4	B	401	5/5	0.18	0.61	25,25,27,32	5
2	PO4	C	401	5/5	0.16	0.10	26,27,27,30	5

6.5 Other polymers ⓘ

There are no such residues in this entry.