



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:09 PM UTC

PDB ID : 2QVS / pdb_00002qvs
Title : Crystal Structure of Type IIa Holoenzyme of cAMP-dependent Protein Kinase
Authors : Wu, J.; Brown, S.H.J.; von Daake, S.; Taylor, S.S.
Deposited on : 2007-08-08
Resolution : 2.50 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) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

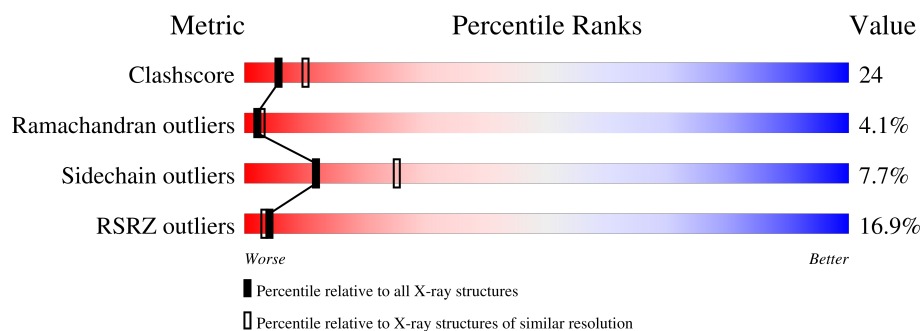
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	E	350	21% 55% 30% 7% 7%
2	B	310	10% 51% 39% 5% . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5056 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 cAMP-dependent protein kinase, alpha-catalytic subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	E	325	Total	C	N	O	P	S	0	0	0
			2525	1632	430	456	2	5			

- Molecule 2 is a protein called cAMP-dependent protein kinase type II-alpha regulatory subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	298	Total	C	N	O	S	0	0	0
			2350	1471	407	457	15			

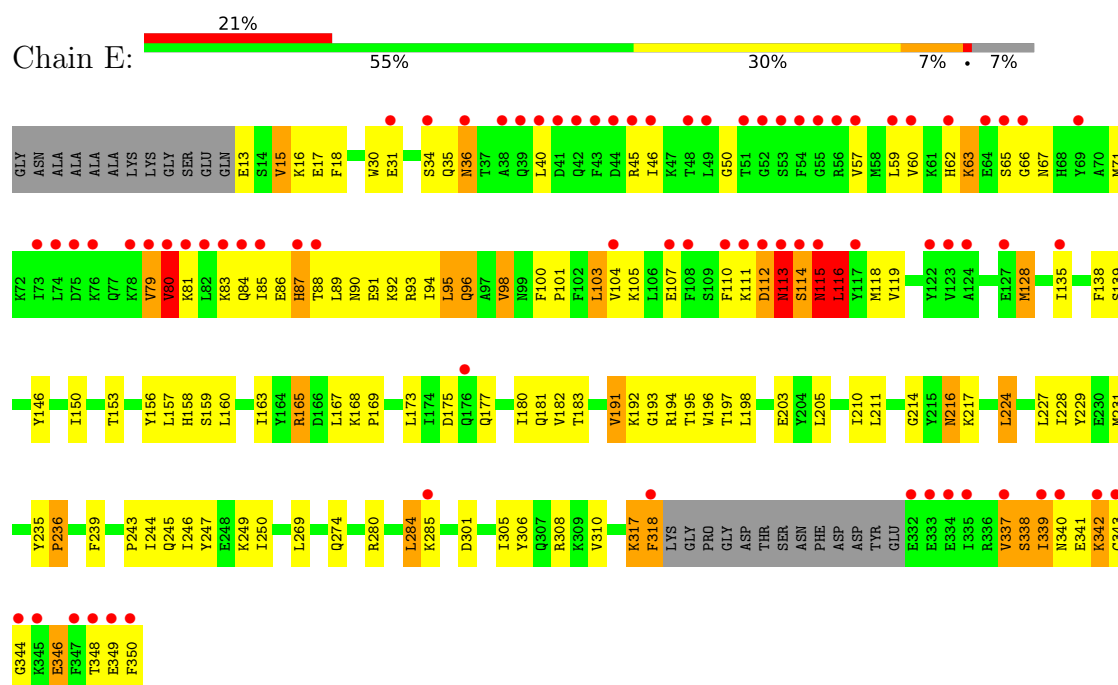
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	106	Total	O	0	0
			106	106		
3	B	75	Total	O	0	0
			75	75		

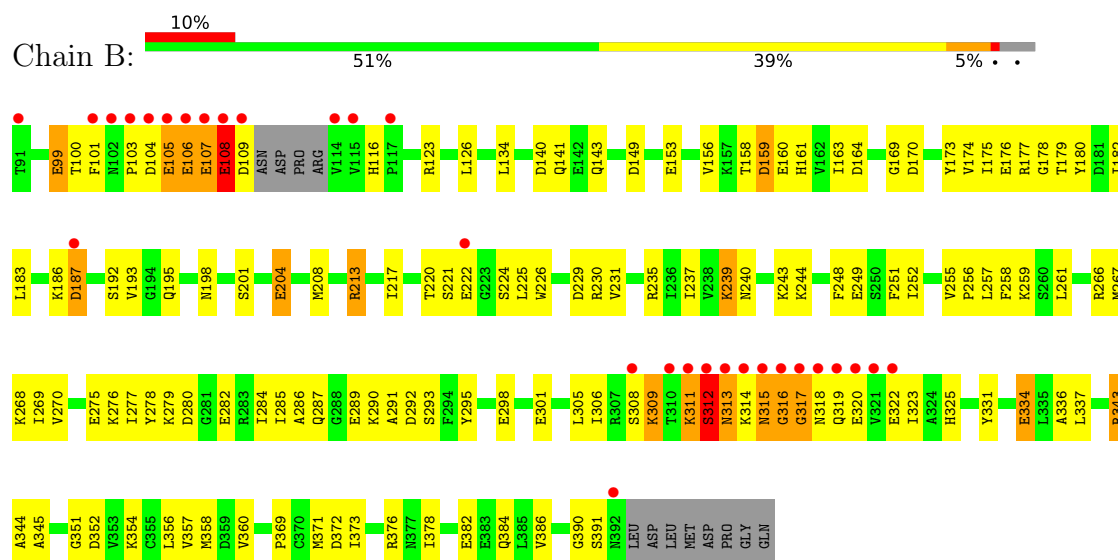
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: cAMP-dependent protein kinase, alpha-catalytic subunit



- Molecule 2: cAMP-dependent protein kinase type II-alpha regulatory subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.90Å 92.90Å 118.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	78.4 (50.00-2.50) 92.1 (50.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.71 (at 2.48Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.212 , 0.251 0.236 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	45.8	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5056	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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

5.1 Standard geometry

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

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	E	0.46	1/2566 (0.0%)	1.17	23/3474 (0.7%)
2	B	0.53	1/2383 (0.0%)	1.06	17/3203 (0.5%)
All	All	0.50	2/4949 (0.0%)	1.12	40/6677 (0.6%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	81	LYS	CA-C	-6.64	1.50	1.53
2	B	311	LYS	CA-CB	5.43	1.60	1.53

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	81	LYS	CA-C-O	22.02	130.71	117.94
1	E	114	SER	CA-C-O	17.25	139.01	120.55
1	E	116	LEU	CA-C-O	15.93	143.29	120.51
1	E	115	ASN	N-CA-C	11.28	134.81	110.80
1	E	114	SER	N-CA-C	-10.60	98.47	111.40
1	E	80	VAL	CB-CA-C	-10.33	94.35	111.29
2	B	314	LYS	N-CA-C	9.80	131.68	110.80
1	E	81	LYS	N-CA-CB	9.74	126.96	110.49
1	E	114	SER	CB-CA-C	8.71	125.42	110.79
1	E	113	ASN	CB-CA-C	-7.55	95.39	110.42
1	E	81	LYS	N-CA-C	-7.50	102.63	108.78
2	B	316	GLY	N-CA-C	7.34	126.05	112.77
2	B	315	ASN	CA-C-O	7.14	130.72	120.51
1	E	114	SER	N-CA-CB	7.02	120.52	110.13
2	B	187	ASP	CB-CA-C	-6.76	106.88	116.34
2	B	309	LYS	N-CA-C	-6.75	104.20	113.18
2	B	314	LYS	CA-C-O	-6.72	110.90	120.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	313	ASN	CA-C-O	-6.58	111.11	120.51
1	E	236	PRO	CA-C-N	6.55	126.49	119.28
1	E	236	PRO	C-N-CA	6.55	126.49	119.28
2	B	116	HIS	CA-C-N	6.52	126.55	119.90
2	B	116	HIS	C-N-CA	6.52	126.55	119.90
1	E	116	LEU	CB-CA-C	-6.49	97.50	110.42
2	B	193	VAL	N-CA-C	6.46	120.10	111.44
2	B	314	LYS	CB-CA-C	-6.41	97.66	110.42
1	E	115	ASN	CB-CA-C	-6.28	97.92	110.42
2	B	378	ILE	N-CA-C	5.77	118.26	111.05
2	B	237	ILE	N-CA-C	5.75	115.94	110.53
1	E	98	VAL	N-CA-C	5.68	117.77	108.97
2	B	134	LEU	N-CA-C	5.67	118.19	111.33
1	E	192	LYS	N-CA-C	-5.63	98.03	107.99
1	E	139	SER	N-CA-C	-5.58	103.16	110.53
2	B	316	GLY	CA-C-O	-5.51	114.17	122.71
2	B	312	SER	N-CA-C	5.43	122.37	110.80
1	E	105	LYS	N-CA-C	5.40	118.21	109.40
1	E	235	TYR	CA-C-N	5.36	123.52	119.66
1	E	235	TYR	C-N-CA	5.36	123.52	119.66
2	B	222	GLU	N-CA-C	-5.26	102.07	109.96
1	E	165	ARG	N-CA-C	5.17	119.54	113.23
1	E	214	GLY	N-CA-C	-5.11	105.34	112.18

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	E	2525	0	2346	119	0
2	B	2350	0	2301	109	0
3	B	75	0	0	6	0
3	E	106	0	0	4	0
All	All	5056	0	4647	224	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 24.

All (224) 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)
2:B:108:GLU:N	2:B:108:GLU:OE2	1.95	0.99
1:E:15:VAL:HG23	1:E:16:LYS:HD2	1.46	0.96
2:B:175:ILE:HA	2:B:225:LEU:HD23	1.54	0.88
1:E:285:LYS:HG2	3:E:363:HOH:O	1.74	0.87
2:B:108:GLU:OE2	2:B:108:GLU:CA	2.22	0.86
1:E:59:LEU:HD12	1:E:60:VAL:H	1.42	0.85
1:E:104:VAL:HG11	1:E:183:THR:HG22	1.57	0.84
1:E:16:LYS:HD2	1:E:16:LYS:H	1.48	0.78
2:B:176:GLU:HG3	2:B:177:ARG:HG3	1.68	0.76
1:E:153:THR:O	1:E:157:LEU:HD13	1.85	0.75
2:B:105:GLU:HG2	2:B:105:GLU:O	1.87	0.74
1:E:112:ASP:O	1:E:114:SER:N	2.21	0.74
2:B:305:LEU:O	2:B:345:ALA:HB1	1.88	0.73
1:E:13:GLU:HB2	1:E:15:VAL:HG22	1.70	0.72
1:E:112:ASP:C	1:E:114:SER:H	1.99	0.71
1:E:86:GLU:C	1:E:88:THR:H	2.00	0.70
1:E:115:ASN:O	1:E:116:LEU:CB	2.40	0.69
1:E:80:VAL:H	1:E:85:ILE:HG21	1.57	0.69
1:E:100:PHE:CD1	1:E:101:PRO:HD2	2.28	0.69
1:E:135:ILE:HD11	1:E:138:PHE:CE1	2.29	0.68
1:E:62:HIS:O	1:E:66:GLY:HA2	1.94	0.68
1:E:113:ASN:HA	1:E:341:GLU:HA	1.74	0.68
2:B:280:ASP:HB2	2:B:351:GLY:HA2	1.75	0.68
1:E:227:LEU:HD13	1:E:231:MET:HE2	1.76	0.67
1:E:337:VAL:HG23	1:E:339:ILE:H	1.60	0.66
2:B:240:ASN:HB2	3:B:449:HOH:O	1.95	0.66
1:E:62:HIS:HB3	1:E:67:ASN:H	1.61	0.65
2:B:140:ASP:OD1	2:B:143:GLN:HG3	1.95	0.65
1:E:40:LEU:HB2	1:E:110:PHE:CE2	2.31	0.65
2:B:235:ARG:NH2	2:B:235:ARG:HB3	2.11	0.65
1:E:35:GLN:HB3	1:E:350:PHE:OXT	1.98	0.64
2:B:107:GLU:O	2:B:108:GLU:C	2.40	0.64
2:B:182:ILE:HG23	2:B:217:ILE:CD1	2.27	0.64
1:E:80:VAL:H	1:E:85:ILE:CG2	2.10	0.64
2:B:108:GLU:OE2	2:B:108:GLU:HA	1.98	0.63
2:B:311:LYS:O	2:B:312:SER:CB	2.46	0.63
1:E:71:MET:HA	1:E:118:MET:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:291:ALA:HB1	2:B:336:ALA:HB2	1.81	0.62
1:E:85:ILE:C	1:E:85:ILE:HD12	2.24	0.62
1:E:146:TYR:HB3	1:E:180:ILE:HD11	1.80	0.62
2:B:331:TYR:CE1	2:B:334:GLU:HG2	2.33	0.62
2:B:256:PRO:HA	2:B:259:LYS:HG2	1.80	0.62
1:E:13:GLU:HB2	1:E:15:VAL:CG2	2.29	0.62
1:E:104:VAL:CG1	1:E:183:THR:HG22	2.28	0.62
2:B:306:ILE:HB	2:B:323:ILE:HD11	1.82	0.62
1:E:128:MET:HA	1:E:128:MET:HE3	1.81	0.61
1:E:227:LEU:CD1	1:E:231:MET:HE2	2.31	0.60
1:E:100:PHE:CG	1:E:101:PRO:HD2	2.35	0.60
1:E:16:LYS:HD2	1:E:16:LYS:N	2.17	0.60
1:E:193:GLY:HA2	2:B:267:MET:SD	2.42	0.60
2:B:235:ARG:CB	2:B:235:ARG:HH21	2.15	0.60
2:B:343:ARG:HG3	2:B:343:ARG:HH11	1.65	0.60
2:B:235:ARG:HB3	2:B:235:ARG:HH21	1.66	0.59
1:E:107:GLU:HB2	1:E:119:VAL:O	2.02	0.59
1:E:342:LYS:O	1:E:344:GLY:N	2.34	0.59
1:E:156:TYR:CD2	1:E:157:LEU:HD12	2.38	0.58
1:E:113:ASN:CA	1:E:341:GLU:HA	2.32	0.58
2:B:101:PHE:HZ	2:B:239:LYS:HZ3	1.50	0.58
1:E:342:LYS:C	1:E:344:GLY:H	2.11	0.58
2:B:176:GLU:HB3	2:B:224:SER:O	2.04	0.58
2:B:266:ARG:O	2:B:270:VAL:HG23	2.04	0.58
1:E:156:TYR:O	1:E:159:SER:HB3	2.04	0.58
1:E:243:PRO:HA	1:E:246:ILE:HD12	1.85	0.58
1:E:62:HIS:CD2	1:E:63:LYS:H	2.21	0.58
1:E:150:ILE:HD13	1:E:167:LEU:HD11	1.86	0.58
2:B:275:GLU:OE2	2:B:354:LYS:HE2	2.03	0.58
2:B:278:TYR:CD2	2:B:284:ILE:HG12	2.39	0.58
2:B:126:LEU:HD23	2:B:174:VAL:HG11	1.86	0.57
2:B:108:GLU:O	2:B:109:ASP:OD1	2.22	0.57
2:B:163:ILE:HB	2:B:213:ARG:HG2	1.86	0.57
1:E:163:ILE:HD12	1:E:165:ARG:HD3	1.85	0.57
1:E:35:GLN:O	1:E:36:ASN:C	2.46	0.57
2:B:182:ILE:C	2:B:183:LEU:HD12	2.30	0.57
2:B:276:LYS:HZ3	2:B:357:VAL:HG21	1.70	0.56
1:E:135:ILE:HD11	1:E:138:PHE:HE1	1.69	0.56
2:B:100:THR:HG22	2:B:101:PHE:N	2.21	0.56
1:E:305:ILE:HD12	1:E:310:VAL:HG21	1.87	0.56
2:B:305:LEU:HD23	2:B:322:GLU:HA	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:107:GLU:OE1	1:E:107:GLU:HA	2.06	0.56
2:B:158:THR:O	2:B:159:ASP:HB2	2.05	0.56
1:E:89:LEU:HD11	1:E:349:GLU:O	2.05	0.56
1:E:113:ASN:N	1:E:341:GLU:HG3	2.22	0.56
2:B:123:ARG:NH1	2:B:149:ASP:OD2	2.39	0.55
2:B:156:VAL:HG13	2:B:160:GLU:OE1	2.07	0.55
1:E:86:GLU:C	1:E:88:THR:N	2.60	0.55
1:E:224:LEU:HD22	1:E:228:ILE:HG13	1.88	0.55
1:E:210:ILE:HG23	2:B:208:MET:HE1	1.89	0.55
2:B:301:GLU:OE1	2:B:325:HIS:HE1	1.90	0.55
1:E:318:PHE:HD1	1:E:318:PHE:O	1.90	0.55
2:B:251:PHE:O	2:B:255:VAL:HG23	2.07	0.54
1:E:40:LEU:HD12	1:E:40:LEU:O	2.08	0.54
2:B:285:ILE:CG2	2:B:343:ARG:HG2	2.37	0.54
2:B:99:GLU:HG3	2:B:231:VAL:HG22	1.90	0.54
1:E:50:GLY:HA3	1:E:57:VAL:O	2.08	0.53
2:B:373:ILE:HA	2:B:376:ARG:HG3	1.89	0.53
2:B:316:GLY:O	2:B:317:GLY:C	2.51	0.53
1:E:18:PHE:HE2	1:E:306:TYR:HH	1.56	0.53
2:B:386:VAL:O	2:B:390:GLY:HA3	2.08	0.53
2:B:309:LYS:C	2:B:311:LYS:H	2.17	0.52
1:E:245:GLN:O	1:E:249:LYS:HG3	2.10	0.52
2:B:276:LYS:HD2	2:B:278:TYR:OH	2.08	0.52
2:B:293:SER:HB2	2:B:358:MET:O	2.09	0.52
1:E:34:SER:HB3	3:E:389:HOH:O	2.09	0.52
1:E:112:ASP:C	1:E:114:SER:N	2.65	0.52
2:B:108:GLU:O	2:B:109:ASP:CG	2.53	0.52
2:B:279:LYS:O	2:B:282:GLU:HG2	2.10	0.52
1:E:15:VAL:HG23	1:E:16:LYS:CD	2.29	0.52
1:E:227:LEU:O	1:E:231:MET:HG3	2.09	0.51
2:B:153:GLU:HB2	2:B:226:TRP:CZ3	2.45	0.51
1:E:59:LEU:HD12	1:E:60:VAL:N	2.18	0.51
2:B:316:GLY:O	2:B:318:ASN:N	2.43	0.51
2:B:286:ALA:O	2:B:289:GLU:HG2	2.11	0.51
1:E:103:LEU:HD21	1:E:157:LEU:HD11	1.93	0.51
1:E:203:GLU:HG2	1:E:236:PRO:HG3	1.93	0.51
2:B:106:GLU:O	2:B:107:GLU:HB2	2.09	0.51
2:B:248:PHE:CE2	2:B:356:LEU:HD21	2.45	0.51
1:E:92:LYS:HD2	1:E:92:LYS:O	2.11	0.50
2:B:169:GLY:HA2	2:B:213:ARG:HH11	1.76	0.50
1:E:191:VAL:HG13	1:E:216:ASN:HD21	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:246:ILE:O	1:E:250:ILE:HG13	2.12	0.50
1:E:236:PRO:HB2	3:E:364:HOH:O	2.12	0.50
2:B:108:GLU:H	2:B:108:GLU:CD	2.18	0.49
2:B:175:ILE:HG12	2:B:225:LEU:HD21	1.94	0.49
2:B:248:PHE:O	2:B:252:ILE:HG12	2.13	0.49
1:E:135:ILE:HD11	1:E:138:PHE:CD1	2.47	0.49
2:B:235:ARG:O	2:B:239:LYS:HB2	2.12	0.49
2:B:107:GLU:O	2:B:109:ASP:OD2	2.30	0.49
1:E:338:SEP:C	1:E:340:ASN:H	2.25	0.49
1:E:111:LYS:CB	1:E:116:LEU:N	2.76	0.49
2:B:244:LYS:HD2	3:B:465:HOH:O	2.13	0.49
2:B:105:GLU:OE1	2:B:243:LYS:HE2	2.13	0.49
2:B:295:TYR:N	2:B:295:TYR:CD2	2.81	0.49
2:B:158:THR:HB	2:B:221:SER:HA	1.94	0.48
1:E:79:VAL:O	1:E:80:VAL:CB	2.61	0.48
1:E:84:GLN:HG3	1:E:85:ILE:N	2.28	0.48
1:E:90:ASN:C	1:E:90:ASN:HD22	2.19	0.48
1:E:111:LYS:CB	1:E:116:LEU:H	2.26	0.48
2:B:268:LYS:HE3	3:B:463:HOH:O	2.12	0.48
2:B:169:GLY:HA2	2:B:213:ARG:NH1	2.29	0.48
1:E:198:LEU:HD11	2:B:208:MET:HE2	1.96	0.48
2:B:280:ASP:HB2	2:B:351:GLY:CA	2.42	0.48
2:B:269:ILE:HD12	2:B:358:MET:HE2	1.96	0.47
2:B:309:LYS:O	2:B:311:LYS:N	2.37	0.47
1:E:84:GLN:HG3	1:E:86:GLU:H	1.79	0.47
1:E:175:ASP:OD2	1:E:308:ARG:NH2	2.45	0.47
2:B:179:THR:C	2:B:180:TYR:CD1	2.93	0.47
1:E:93:ARG:O	1:E:96:GLN:HG3	2.15	0.47
1:E:338:SEP:C	1:E:340:ASN:N	2.74	0.47
2:B:175:ILE:HG12	2:B:225:LEU:CD2	2.44	0.47
2:B:285:ILE:HG21	2:B:343:ARG:HG2	1.95	0.47
2:B:183:LEU:HD12	2:B:183:LEU:N	2.30	0.47
2:B:105:GLU:O	2:B:105:GLU:CG	2.62	0.47
1:E:104:VAL:HG13	1:E:182:VAL:O	2.15	0.46
2:B:258:PHE:HB3	2:B:261:LEU:HD12	1.97	0.46
1:E:301:ASP:O	1:E:305:ILE:HD13	2.14	0.46
1:E:158:HIS:HB3	1:E:217:LYS:HD3	1.96	0.46
2:B:309:LYS:C	2:B:311:LYS:N	2.73	0.46
1:E:86:GLU:O	1:E:88:THR:N	2.48	0.46
1:E:317:LYS:HB2	1:E:317:LYS:NZ	2.31	0.46
2:B:106:GLU:H	2:B:106:GLU:HG3	1.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:343:ARG:HG3	2:B:343:ARG:NH1	2.30	0.46
2:B:278:TYR:HA	2:B:282:GLU:OE2	2.16	0.46
2:B:280:ASP:HA	2:B:351:GLY:O	2.15	0.46
1:E:156:TYR:HD2	1:E:157:LEU:HD12	1.80	0.46
2:B:249:GLU:OE2	2:B:249:GLU:HA	2.16	0.46
1:E:35:GLN:HA	1:E:350:PHE:O	2.16	0.45
1:E:177:GLN:NE2	1:E:308:ARG:NH1	2.64	0.45
1:E:191:VAL:HG13	1:E:216:ASN:ND2	2.29	0.45
1:E:244:ILE:HD13	2:B:204:GLU:OE1	2.16	0.45
2:B:161:HIS:ND1	2:B:164:ASP:OD1	2.46	0.45
2:B:293:SER:HA	2:B:360:VAL:HG23	1.98	0.45
1:E:84:GLN:HG3	1:E:85:ILE:H	1.81	0.45
2:B:317:GLY:O	2:B:318:ASN:C	2.58	0.45
1:E:36:ASN:HA	1:E:110:PHE:HA	1.98	0.45
2:B:107:GLU:O	2:B:109:ASP:CG	2.60	0.45
2:B:298:GLU:HG2	2:B:354:LYS:O	2.17	0.45
1:E:280:ARG:O	1:E:284:LEU:HD13	2.17	0.44
2:B:295:TYR:O	2:B:331:TYR:HB2	2.17	0.44
1:E:342:LYS:C	1:E:344:GLY:N	2.76	0.44
2:B:178:GLY:N	2:B:198:ASN:OD1	2.50	0.44
1:E:65:SER:CB	1:E:67:ASN:ND2	2.81	0.44
1:E:195:THR:C	1:E:196:TRP:CD1	2.96	0.44
2:B:173:TYR:O	2:B:201:SER:HA	2.18	0.44
1:E:94:ILE:O	1:E:98:VAL:HG22	2.17	0.44
1:E:128:MET:HG2	1:E:169:PRO:HA	2.00	0.44
2:B:306:ILE:O	2:B:306:ILE:HG23	2.18	0.43
1:E:85:ILE:C	1:E:85:ILE:CD1	2.91	0.43
2:B:287:GLN:HG3	2:B:344:ALA:O	2.18	0.43
1:E:83:LYS:O	1:E:83:LYS:HD3	2.19	0.43
1:E:80:VAL:N	1:E:85:ILE:HG21	2.29	0.43
2:B:292:ASP:O	2:B:360:VAL:HG23	2.18	0.43
2:B:158:THR:HB	2:B:221:SER:CA	2.49	0.43
1:E:167:LEU:O	1:E:168:LYS:HB3	2.18	0.43
1:E:346:GLU:C	1:E:348:THR:H	2.27	0.42
2:B:220:THR:HG21	3:B:446:HOH:O	2.19	0.42
1:E:91:GLU:O	1:E:95:LEU:HD23	2.19	0.42
2:B:291:ALA:HB3	3:B:404:HOH:O	2.19	0.42
2:B:306:ILE:CB	2:B:323:ILE:HD11	2.48	0.42
2:B:320:GLU:OE1	2:B:320:GLU:HA	2.20	0.42
1:E:36:ASN:HD22	1:E:111:LYS:H	1.67	0.42
1:E:91:GLU:O	1:E:95:LEU:CD2	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:193:GLY:O	1:E:194:ARG:NH1	2.49	0.42
1:E:341:GLU:CD	1:E:341:GLU:N	2.78	0.42
1:E:92:LYS:HD2	1:E:92:LYS:C	2.45	0.42
2:B:140:ASP:OD2	2:B:143:GLN:NE2	2.52	0.42
1:E:103:LEU:HD12	1:E:182:VAL:HB	2.01	0.42
1:E:156:TYR:HD2	1:E:157:LEU:CD1	2.33	0.42
1:E:13:GLU:N	1:E:16:LYS:HD3	2.35	0.42
1:E:216:ASN:HB3	3:E:402:HOH:O	2.20	0.42
1:E:229:TYR:CD1	1:E:229:TYR:C	2.98	0.41
1:E:203:GLU:OE2	1:E:239:PHE:HA	2.19	0.41
2:B:371:MET:HE3	3:B:456:HOH:O	2.20	0.41
2:B:204:GLU:CD	2:B:204:GLU:H	2.28	0.41
2:B:369:PRO:HG2	2:B:372:ASP:OD2	2.21	0.41
1:E:90:ASN:C	1:E:90:ASN:ND2	2.77	0.41
1:E:173:LEU:HD13	1:E:183:THR:CG2	2.51	0.41
1:E:205:LEU:HD12	1:E:247:TYR:CE1	2.56	0.41
2:B:103:PRO:HG2	2:B:104:ASP:H	1.85	0.41
2:B:170:ASP:HA	2:B:230:ARG:HH12	1.84	0.41
2:B:337:LEU:HD23	2:B:360:VAL:HG13	2.03	0.40
2:B:186:LYS:O	2:B:187:ASP:HB2	2.20	0.40
2:B:276:LYS:C	2:B:277:ILE:HD12	2.45	0.40
1:E:317:LYS:NZ	1:E:317:LYS:CB	2.85	0.40
1:E:30:TRP:HZ3	1:E:94:ILE:HG12	1.87	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	E	319/350 (91%)	275 (86%)	28 (9%)	16 (5%)	1	2
2	B	294/310 (95%)	270 (92%)	15 (5%)	9 (3%)	3	5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	613/660 (93%)	545 (89%)	43 (7%)	25 (4%)	2 3

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	79	VAL
1	E	112	ASP
1	E	115	ASN
1	E	343	CYS
2	B	107	GLU
2	B	308	SER
2	B	312	SER
2	B	315	ASN
2	B	317	GLY
1	E	36	ASN
1	E	80	VAL
1	E	113	ASN
1	E	337	VAL
2	B	391	SER
1	E	46	ILE
1	E	116	LEU
2	B	108	GLU
2	B	313	ASN
2	B	319	GLN
1	E	45	ARG
1	E	63	LYS
1	E	87	HIS
1	E	216	ASN
1	E	346	GLU
1	E	339	ILE

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	E	238/303 (78%)	219 (92%)	19 (8%)	11 24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	253/273 (93%)	234 (92%)	19 (8%)	12	26
All	All	491/576 (85%)	453 (92%)	38 (8%)	12	25

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	15	VAL
1	E	17	GLU
1	E	31	GLU
1	E	87	HIS
1	E	95	LEU
1	E	96	GLN
1	E	103	LEU
1	E	128	MET
1	E	160	LEU
1	E	181	GLN
1	E	191	VAL
1	E	211	LEU
1	E	224	LEU
1	E	269	LEU
1	E	274	GLN
1	E	284	LEU
1	E	317	LYS
1	E	318	PHE
1	E	342	LYS
2	B	99	GLU
2	B	105	GLU
2	B	106	GLU
2	B	108	GLU
2	B	141	GLN
2	B	159	ASP
2	B	192	SER
2	B	195	GLN
2	B	204	GLU
2	B	213	ARG
2	B	229	ASP
2	B	239	LYS
2	B	257	LEU
2	B	290	LYS
2	B	334	GLU
2	B	343	ARG
2	B	352	ASP

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Mol	Chain	Res	Type
2	B	382	GLU
2	B	384	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	35	GLN
1	E	36	ASN
1	E	39	GLN
1	E	62	HIS
1	E	67	ASN
1	E	87	HIS
1	E	90	ASN
1	E	96	GLN
1	E	171	ASN
1	E	176	GLN
1	E	177	GLN
1	E	181	GLN
1	E	216	ASN
1	E	245	GLN
1	E	289	ASN
1	E	293	ASN
2	B	122	GLN
2	B	127	GLN
2	B	141	GLN
2	B	143	GLN
2	B	146	GLN
2	B	171	ASN
2	B	195	GLN
2	B	325	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	TPO	E	197	1	8,10,11	0.93	1 (12%)	10,14,16	1.01	0
1	SEP	E	338	1	8,9,10	0.72	0	7,12,14	1.43	1 (14%)

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
1	TPO	E	197	1	-	1/9/11/13	-
1	SEP	E	338	1	-	3/6/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	197	TPO	P-OG1	2.10	1.63	1.59

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	338	SEP	OG-CB-CA	3.38	111.44	108.14

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	E	338	SEP	N-CA-CB-OG
1	E	338	SEP	C-CA-CB-OG
1	E	338	SEP	CA-CB-OG-P
1	E	197	TPO	CA-CB-OG1-P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	338	SEP	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 ⓘ

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	E	323/350 (92%)	0.95	75 (23%) 2 1	21, 42, 100, 101	0
2	B	298/310 (96%)	0.71	30 (10%) 12 10	27, 43, 79, 91	0
All	All	621/660 (94%)	0.83	105 (16%) 4 3	21, 43, 93, 101	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	315	ASN	7.7
2	B	310	THR	7.3
2	B	114	VAL	6.5
2	B	106	GLU	6.2
2	B	108	GLU	6.1
2	B	312	SER	5.4
2	B	308	SER	5.2
1	E	85	ILE	5.2
2	B	313	ASN	5.1
1	E	54	PHE	5.1
2	B	318	ASN	5.0
1	E	80	VAL	4.6
1	E	45	ARG	4.5
1	E	44	ASP	4.5
1	E	82	LEU	4.4
1	E	40	LEU	4.3
2	B	311	LYS	4.1
1	E	43	PHE	4.0
1	E	318	PHE	4.0
1	E	49	LEU	4.0
1	E	51	THR	3.9
2	B	105	GLU	3.9
1	E	114	SER	3.9
1	E	350	PHE	3.9

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Mol	Chain	Res	Type	RSRZ
2	B	109	ASP	3.9
1	E	53	SER	3.9
1	E	110	PHE	3.8
1	E	52	GLY	3.8
2	B	107	GLU	3.8
1	E	335	ILE	3.8
1	E	83	LYS	3.8
1	E	87	HIS	3.8
1	E	347	PHE	3.7
2	B	317	GLY	3.7
1	E	84	GLN	3.7
2	B	115	VAL	3.6
1	E	108	PHE	3.6
1	E	124	ALA	3.6
2	B	319	GLN	3.6
2	B	103	PRO	3.5
1	E	56	ARG	3.5
1	E	38	ALA	3.5
1	E	334	GLU	3.5
1	E	333	GLU	3.4
1	E	176	GLN	3.3
1	E	339	ILE	3.3
1	E	64	GLU	3.3
1	E	123	VAL	3.3
1	E	117	TYR	3.2
2	B	316	GLY	3.2
1	E	343	CYS	3.2
1	E	42	GLN	3.2
1	E	75	ASP	3.2
1	E	332	GLU	3.1
1	E	46	ILE	3.1
1	E	81	LYS	3.0
1	E	88	THR	2.9
1	E	115	ASN	2.9
1	E	285	LYS	2.9
1	E	337	VAL	2.8
1	E	127	GLU	2.8
2	B	104	ASP	2.7
1	E	76	LYS	2.7
2	B	102	ASN	2.7
1	E	349	GLU	2.7
1	E	57	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	91	THR	2.6
1	E	39	GLN	2.6
1	E	69	TYR	2.6
1	E	31	GLU	2.6
1	E	59	LEU	2.6
1	E	79	VAL	2.6
1	E	34	SER	2.6
1	E	342	LYS	2.6
2	B	222	GLU	2.5
1	E	62	HIS	2.5
1	E	78	LYS	2.5
1	E	48	THR	2.5
2	B	320	GLU	2.5
1	E	55	GLY	2.5
1	E	41	ASP	2.5
2	B	392	ASN	2.5
2	B	321	VAL	2.4
1	E	107	GLU	2.4
1	E	113	ASN	2.3
1	E	73	ILE	2.3
1	E	112	ASP	2.3
1	E	36	ASN	2.3
1	E	74	LEU	2.3
2	B	187	ASP	2.3
1	E	60	VAL	2.2
1	E	340	ASN	2.2
1	E	122	TYR	2.2
1	E	111	LYS	2.2
1	E	348	THR	2.2
1	E	104	VAL	2.1
1	E	65	SER	2.1
2	B	322	GLU	2.1
2	B	314	LYS	2.1
1	E	135	ILE	2.1
1	E	66	GLY	2.0
1	E	344	GLY	2.0
1	E	345	LYS	2.0
2	B	117	PRO	2.0
2	B	101	PHE	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	SEP	E	338	10/11	0.75	0.12	100,100,100,100	0
1	TPO	E	197	11/12	0.97	0.07	28,29,30,32	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.