



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:33 AM UTC

PDB ID : 2WH0 / pdb_00002wh0
Title : Recognition of an intrachain tandem 14-3-3 binding site within protein kinase C epsilon
Authors : Kostecky, B.; Saurin, A.T.; Purkiss, A.; Parker, P.J.; McDonald, N.Q.
Deposited on : 2009-04-28
Resolution : 2.25 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:

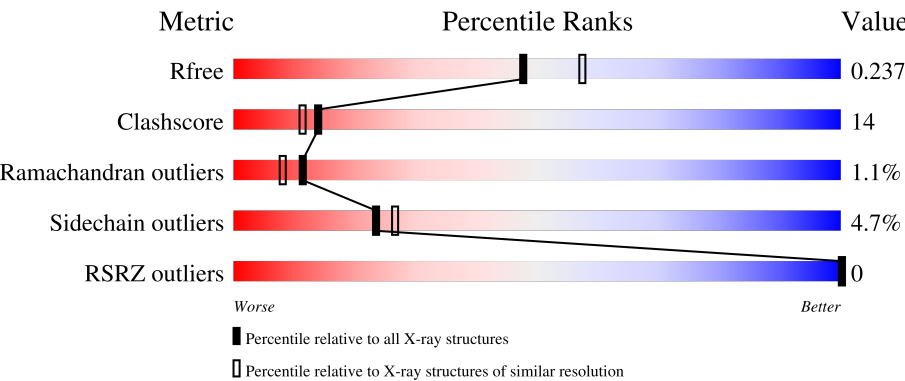
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	180053	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	A	245	63%25%•10%
1	B	245	64%27%•7%
1	C	245	65%24%•8%
1	D	245	62%27%•9%
2	Q	31	23%16%•58%

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Mol	Chain	Length	Quality of chain
2	R	31	 19% 10% 71%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6877 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.

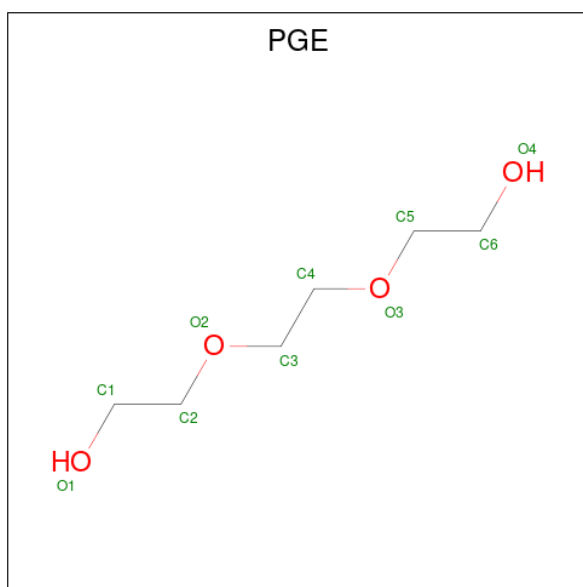
- Molecule 1 is a protein called 14-3-3 PROTEIN ZETA/DELTA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1631	1035	270	317	9			
1	B	227	Total	C	N	O	S	0	0	0
			1741	1100	286	346	9			
1	C	226	Total	C	N	O	S	0	0	0
			1681	1062	285	325	9			
1	D	222	Total	C	N	O	S	0	0	0
			1610	1016	271	314	9			

- Molecule 2 is a protein called PROTEIN KINASE C EPSILON TYPE, NPKC-EPSILON.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Q	13	Total	C	N	O	P	0	0	0
			84	45	14	23	2			
2	R	9	Total	C	N	O	P	0	0	0
			58	30	9	17	2			

- Molecule 3 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Ca	0	0
			1	1		

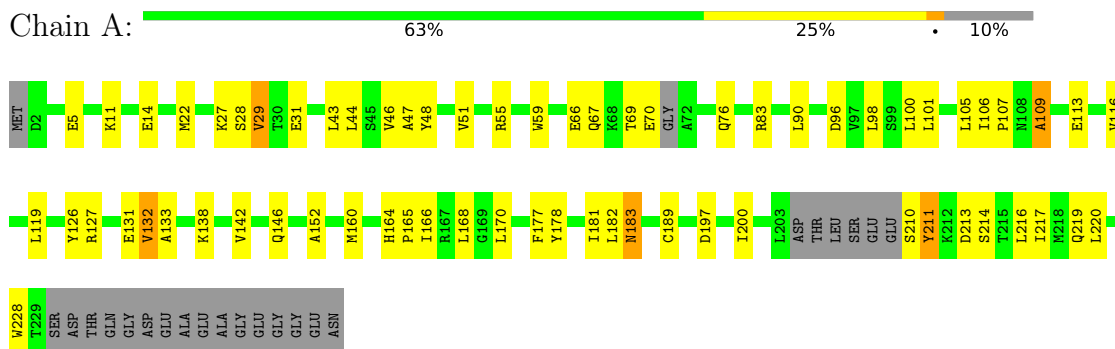
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	15	Total	O	0	0
			15	15		
5	B	24	Total	O	0	0
			24	24		
5	C	14	Total	O	0	0
			14	14		
5	D	8	Total	O	0	0
			8	8		
5	Q	3	Total	O	0	0
			3	3		

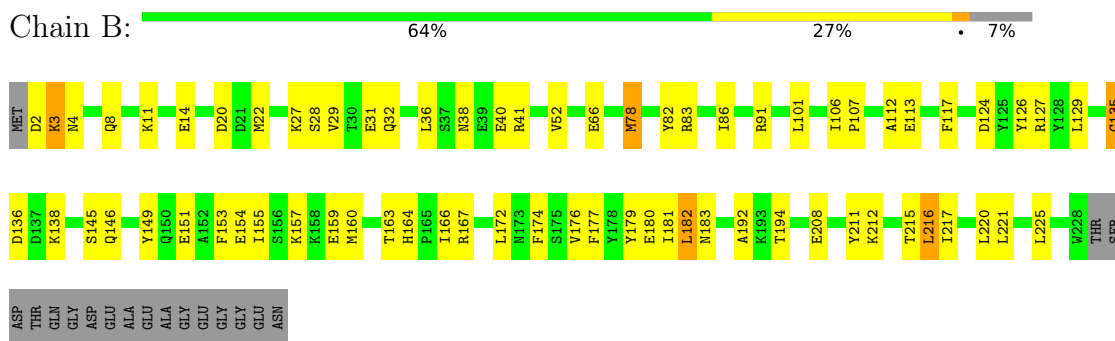
3 Residue-property plots [i](#)

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.

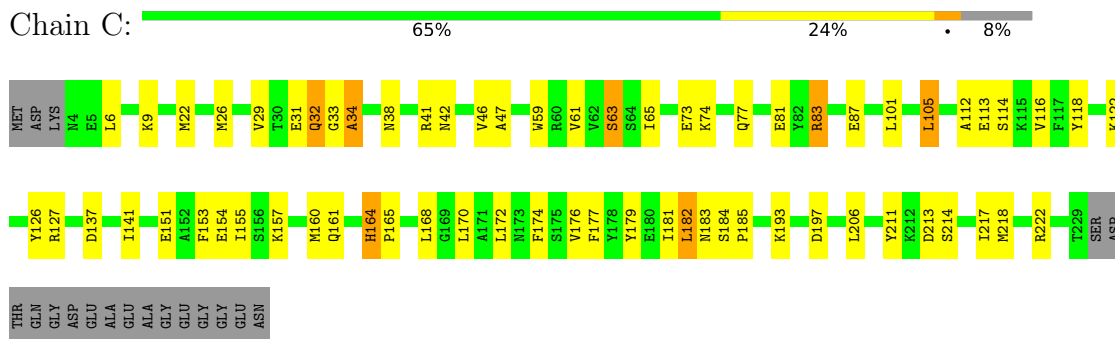
• Molecule 1: 14-3-3 PROTEIN ZETA/DELTA



• Molecule 1: 14-3-3 PROTEIN ZETA/DELTA

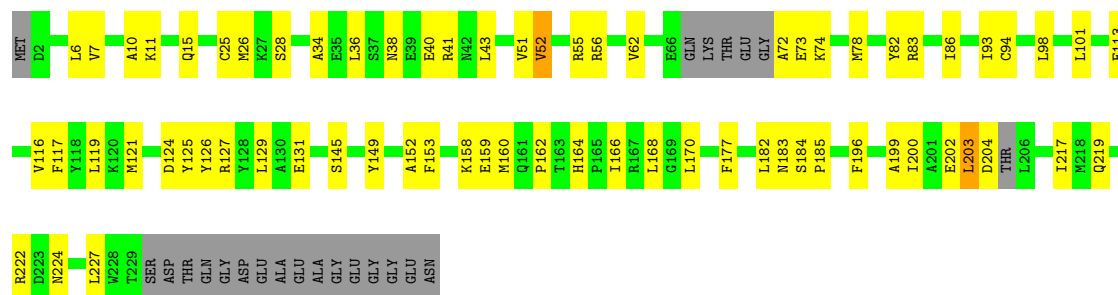


• Molecule 1: 14-3-3 PROTEIN ZETA/DELTA



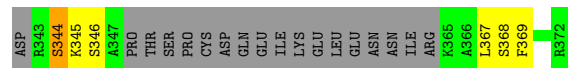
- Molecule 1: 14-3-3 PROTEIN ZETA/DELTA

Chain D:  62% 27% 9%



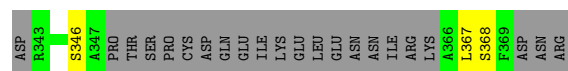
- Molecule 2: PROTEIN KINASE C EPSILON TYPE, NPKC-EPSILON

Chain Q:  23% 16% 58%



- Molecule 2: PROTEIN KINASE C EPSILON TYPE, NPKC-EPSILON

Chain R:  19% 10% 71%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.13Å 78.16Å 108.52Å 90.00° 90.10° 90.00°	Depositor
Resolution (Å)	24.46 – 2.25 24.46 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.4 (24.46-2.25) 99.2 (24.46-2.25)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.48 (at 2.26Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.181 , 0.235 0.185 , 0.237	Depositor DCC
R_{free} test set	1996 reflections (3.55%)	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.560	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 15.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.420 for h,-k,-l	Xtriage
Reported twinning fraction	0.441 for h,-k,-l	Depositor
Outliers	2 of 56277 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6877	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.69% 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, PGE, 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	0.40	0/1653	0.78	0/2241
1	B	0.41	0/1766	0.82	1/2386 (0.0%)
1	C	0.39	0/1705	0.85	3/2307 (0.1%)
1	D	0.38	0/1633	0.85	2/2219 (0.1%)
2	Q	1.28	1/60 (1.7%)	1.03	0/74
2	R	0.23	0/34	0.37	0/41
All	All	0.41	1/6851 (0.0%)	0.83	6/9268 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	344	SER	C-O	-5.84	1.16	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	184	SER	CA-C-N	6.70	126.49	119.05
1	D	184	SER	C-N-CA	6.70	126.49	119.05
1	C	164	HIS	CA-C-N	5.54	125.90	119.47
1	C	164	HIS	C-N-CA	5.54	125.90	119.47
1	B	20	ASP	N-CA-C	-5.14	105.76	111.36
1	C	105	LEU	N-CA-C	5.03	117.42	111.33

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	1631	0	1493	51	0
1	B	1741	0	1649	48	0
1	C	1681	0	1569	37	0
1	D	1610	0	1422	53	0
2	Q	84	0	55	4	0
2	R	58	0	34	3	0
3	B	7	0	9	2	0
4	D	1	0	0	0	0
5	A	15	0	0	1	0
5	B	24	0	0	2	0
5	C	14	0	0	0	0
5	D	8	0	0	0	0
5	Q	3	0	0	0	0
All	All	6877	0	6231	188	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 14.

All (188) 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:D:113:GLU:HA	1:D:160:MET:HE2	1.34	1.10
1:D:168:LEU:HB3	1:D:217:ILE:HG21	1.51	0.90
1:D:224:ASN:HD21	2:R:367:LEU:H	1.21	0.88
1:C:101:LEU:HA	1:C:105:LEU:HB2	1.56	0.87
1:D:158:LYS:HG3	1:D:159:GLU:HG2	1.62	0.82
1:D:113:GLU:HA	1:D:160:MET:CE	2.10	0.81
1:A:164:HIS:HE1	1:A:166:ILE:HD12	1.46	0.79
1:B:167:ARG:HG3	3:B:1229:PGE:H42	1.65	0.76
1:D:126:TYR:HB3	1:D:145:SER:HB2	1.69	0.75
1:A:69:THR:HG22	1:A:76:GLN:HB2	1.68	0.74
1:D:119:LEU:HB2	1:D:152:ALA:HB2	1.72	0.71
1:D:183:ASN:O	1:D:185:PRO:HD3	1.91	0.69
1:A:168:LEU:HB3	1:A:217:ILE:HG21	1.73	0.69
1:D:224:ASN:ND2	2:R:367:LEU:H	1.89	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:113:GLU:CA	1:D:160:MET:HE2	2.20	0.68
1:A:220:LEU:HD22	2:Q:345:LYS:HD3	1.76	0.67
1:D:168:LEU:HD21	1:D:199:ALA:HB2	1.76	0.67
1:B:146:GLN:HG3	1:B:177:PHE:CE2	2.32	0.65
1:A:168:LEU:HD12	1:A:214:SER:HB2	1.79	0.65
1:C:183:ASN:C	1:C:185:PRO:HD3	2.22	0.64
1:B:153:PHE:CE2	1:B:157:LYS:HE2	2.33	0.63
1:B:164:HIS:CE1	1:B:166:ILE:HG12	2.33	0.63
1:A:214:SER:HA	1:A:217:ILE:HD12	1.80	0.63
1:C:31:GLU:O	1:C:32:GLN:HG3	1.99	0.63
1:C:214:SER:O	1:C:218:MET:HG3	1.99	0.62
1:A:5:GLU:O	1:B:78:MET:HE1	2.00	0.62
1:B:220:LEU:HD23	2:Q:367:LEU:HD13	1.82	0.61
1:D:196:PHE:CE1	1:D:222:ARG:HB2	2.34	0.61
1:A:70:GLU:HA	1:A:76:GLN:OE1	2.01	0.60
1:A:116:VAL:HB	1:A:160:MET:HE1	1.83	0.60
1:A:98:LEU:HD23	1:A:101:LEU:HD12	1.82	0.60
1:D:11:LYS:O	1:D:15:GLN:HG3	2.02	0.60
1:B:177:PHE:O	1:B:181:ILE:HB	2.02	0.60
1:A:133:ALA:O	1:A:138:LYS:HB2	2.02	0.59
1:C:22:MET:HG2	1:C:47:ALA:HB2	1.83	0.59
1:B:212:LYS:O	1:B:216:LEU:HB2	2.03	0.58
1:B:192:ALA:HB3	1:B:225:LEU:HD21	1.85	0.58
1:D:116:VAL:HG21	1:D:160:MET:HE3	1.85	0.57
1:D:153:PHE:HA	1:D:170:LEU:HD21	1.86	0.57
1:B:164:HIS:HE1	1:B:166:ILE:HG12	1.70	0.57
1:A:119:LEU:HB3	1:A:152:ALA:HB2	1.86	0.56
1:C:113:GLU:HA	1:C:160:MET:CE	2.35	0.56
1:A:55:ARG:HB3	1:A:90:LEU:HD12	1.86	0.56
1:B:8:GLN:OE1	1:B:11:LYS:HE2	2.06	0.56
1:A:164:HIS:CE1	1:A:166:ILE:HD12	2.34	0.56
1:B:4:ASN:O	1:B:8:GLN:HG2	2.05	0.56
1:A:47:ALA:O	1:A:51:VAL:HG23	2.06	0.56
1:B:179:TYR:O	1:B:183:ASN:HA	2.06	0.56
1:A:213:ASP:O	1:A:217:ILE:HG13	2.07	0.55
1:A:127:ARG:HG3	1:A:181:ILE:HG13	1.89	0.55
1:C:179:TYR:CE2	1:C:185:PRO:HB3	2.41	0.55
1:C:83:ARG:O	1:C:87:GLU:HB2	2.07	0.55
1:B:2:ASP:O	1:B:3:LYS:C	2.49	0.54
1:C:153:PHE:CE2	1:C:157:LYS:HD2	2.43	0.54
1:B:106:ILE:HB	1:B:107:PRO:HD3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:ASP:O	1:B:127:ARG:HB3	2.07	0.53
1:D:98:LEU:HA	1:D:101:LEU:HD12	1.90	0.53
1:D:158:LYS:HG3	1:D:159:GLU:N	2.21	0.53
1:B:146:GLN:HG3	1:B:177:PHE:HE2	1.73	0.53
1:A:146:GLN:OE1	1:A:182:LEU:HD11	2.09	0.52
1:B:220:LEU:HD22	2:Q:367:LEU:HB3	1.90	0.52
1:A:127:ARG:O	1:A:131:GLU:HG3	2.10	0.52
1:A:216:LEU:O	1:A:219:GLN:HB3	2.10	0.52
1:C:213:ASP:O	1:C:217:ILE:HG13	2.10	0.52
1:A:113:GLU:HA	1:A:160:MET:HE2	1.92	0.52
1:D:7:VAL:HG21	1:D:40:GLU:HG2	1.91	0.52
1:A:69:THR:HG22	1:A:69:THR:O	2.08	0.52
1:C:168:LEU:HB3	1:C:217:ILE:HG21	1.92	0.52
1:A:197:ASP:O	1:A:200:ILE:HG22	2.10	0.51
1:C:206:LEU:HD21	1:C:214:SER:CB	2.40	0.51
1:A:22:MET:HG2	1:A:47:ALA:HB2	1.93	0.51
1:C:33:GLY:O	1:C:34:ALA:C	2.54	0.51
1:D:183:ASN:C	1:D:185:PRO:HD3	2.36	0.51
1:D:224:ASN:HD21	2:R:367:LEU:N	2.01	0.50
1:B:41:ARG:HD3	5:B:2003:HOH:O	2.11	0.50
1:B:27:LYS:O	1:B:31:GLU:HG3	2.12	0.50
1:A:27:LYS:HD2	1:A:48:TYR:OH	2.11	0.50
1:B:8:GLN:OE1	1:B:8:GLN:HA	2.10	0.50
1:C:61:VAL:O	1:C:65:ILE:HG13	2.12	0.50
1:D:116:VAL:HG13	1:D:152:ALA:HB1	1.92	0.50
1:D:127:ARG:O	1:D:131:GLU:HG3	2.12	0.50
1:C:6:LEU:HD23	1:C:9:LYS:HD2	1.93	0.50
1:D:38:ASN:HD22	1:D:38:ASN:N	2.08	0.50
1:C:206:LEU:HD22	1:C:211:TYR:HA	1.94	0.50
1:C:41:ARG:HG3	1:C:118:TYR:OH	2.12	0.49
1:A:132:VAL:HG23	5:A:2011:HOH:O	2.12	0.49
1:B:211:TYR:CE2	1:B:215:THR:HG21	2.47	0.49
1:C:63:SER:HB2	1:C:83:ARG:HD3	1.94	0.49
1:D:82:TYR:O	1:D:86:ILE:HG13	2.12	0.49
1:A:210:SER:O	1:A:211:TYR:C	2.55	0.49
1:C:177:PHE:CE2	1:C:182:LEU:HD13	2.47	0.49
1:B:221:LEU:O	1:B:225:LEU:HG	2.13	0.49
1:C:153:PHE:O	1:C:157:LYS:HG3	2.13	0.48
1:D:52:VAL:O	1:D:56:ARG:HG3	2.12	0.48
1:B:113:GLU:HA	1:B:160:MET:CE	2.43	0.48
1:C:151:GLU:O	1:C:155:ILE:HG13	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:42:ASN:O	1:C:46:VAL:HG23	2.13	0.48
1:D:126:TYR:CB	1:D:145:SER:HB2	2.42	0.48
1:D:116:VAL:CG2	1:D:160:MET:HE3	2.44	0.47
1:B:29:VAL:HG12	1:B:36:LEU:HD21	1.96	0.47
1:C:127:ARG:HG3	1:C:181:ILE:HG13	1.95	0.47
1:B:126:TYR:HB3	1:B:145:SER:HB2	1.96	0.47
1:A:66:GLU:OE1	1:A:83:ARG:HD3	2.15	0.47
1:C:26:MET:HA	1:C:29:VAL:HB	1.97	0.47
1:D:10:ALA:HB2	1:D:25:CYS:CB	2.44	0.47
1:D:119:LEU:CB	1:D:152:ALA:HB2	2.42	0.47
1:D:127:ARG:HD3	1:D:177:PHE:HB2	1.95	0.47
1:D:200:ILE:HG23	1:D:203:LEU:HD12	1.97	0.47
1:C:59:TRP:O	1:C:63:SER:HB3	2.15	0.47
1:A:69:THR:O	1:A:69:THR:CG2	2.63	0.47
1:B:160:MET:O	1:B:167:ARG:NH1	2.44	0.47
1:B:112:ALA:HB1	1:B:159:GLU:HG3	1.97	0.46
1:A:116:VAL:CB	1:A:160:MET:HE1	2.45	0.46
1:D:72:ALA:O	1:D:74:LYS:N	2.49	0.46
1:A:165:PRO:HG3	1:A:214:SER:HB3	1.98	0.46
1:D:7:VAL:HG22	1:D:26:MET:HE1	1.98	0.46
1:A:119:LEU:CB	1:A:152:ALA:HB2	2.46	0.46
1:A:168:LEU:CB	1:A:217:ILE:HG21	2.45	0.46
1:B:38:ASN:ND2	5:B:2003:HOH:O	2.49	0.46
1:A:182:LEU:HD23	1:A:182:LEU:HA	1.73	0.45
1:B:182:LEU:O	1:B:183:ASN:C	2.59	0.45
1:B:41:ARG:HD2	1:B:117:PHE:CD1	2.52	0.45
1:B:151:GLU:O	1:B:155:ILE:HG13	2.16	0.45
1:A:113:GLU:HA	1:A:160:MET:CE	2.47	0.45
1:A:142:VAL:HG13	1:A:177:PHE:CE1	2.52	0.45
1:B:172:LEU:O	1:B:176:VAL:HG23	2.17	0.45
1:C:38:ASN:O	1:C:42:ASN:ND2	2.50	0.45
1:D:62:VAL:HG12	1:D:83:ARG:HB2	1.98	0.45
1:B:217:ILE:HA	1:B:220:LEU:HD12	1.98	0.44
1:A:44:LEU:HD21	1:A:105:LEU:HD11	1.98	0.44
1:A:168:LEU:HB3	1:A:217:ILE:CG2	2.43	0.44
1:A:146:GLN:HB2	1:A:177:PHE:CZ	2.53	0.44
1:D:55:ARG:N	1:D:55:ARG:HD2	2.31	0.44
1:B:82:TYR:O	1:B:86:ILE:HG13	2.17	0.44
1:B:28:SER:O	1:B:32:GLN:HG3	2.18	0.44
1:B:113:GLU:HA	1:B:160:MET:HE3	2.00	0.44
1:D:124:ASP:O	1:D:127:ARG:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:55:ARG:HD3	1:D:93:ILE:CD1	2.47	0.43
1:A:106:ILE:N	1:A:107:PRO:CD	2.81	0.43
1:C:77:GLN:O	1:C:81:GLU:HG3	2.18	0.43
1:D:10:ALA:HB2	1:D:25:CYS:HB2	2.01	0.43
1:D:164:HIS:HE1	1:D:166:ILE:HD12	1.83	0.43
1:C:193:LYS:HG2	1:C:197:ASP:OD2	2.19	0.43
1:D:6:LEU:HD13	1:D:28:SER:HB3	2.00	0.43
1:D:7:VAL:HG13	1:D:43:LEU:CD1	2.48	0.43
1:B:36:LEU:HD22	1:B:40:GLU:HB3	2.01	0.43
1:C:33:GLY:O	1:C:34:ALA:O	2.36	0.43
1:A:67:GLN:C	1:A:69:THR:H	2.26	0.43
1:B:135:GLY:O	1:B:138:LYS:HB3	2.18	0.43
1:C:184:SER:N	1:C:185:PRO:HD3	2.33	0.43
1:C:172:LEU:O	1:C:176:VAL:HG23	2.19	0.43
1:B:225:LEU:HA	1:B:225:LEU:HD23	1.72	0.43
1:D:41:ARG:HD3	1:D:117:PHE:CE1	2.53	0.43
1:A:96:ASP:O	1:A:100:LEU:HG	2.18	0.43
1:D:121:MET:O	1:D:125:TYR:HD1	2.02	0.42
1:A:27:LYS:O	1:A:31:GLU:HG3	2.18	0.42
1:C:83:ARG:HD2	1:C:87:GLU:OE1	2.19	0.42
1:B:14:GLU:HB2	1:B:22:MET:CE	2.50	0.42
1:C:137:ASP:O	1:C:141:ILE:HG12	2.20	0.42
1:A:106:ILE:O	1:A:109:ALA:HB3	2.19	0.42
1:A:14:GLU:OE2	1:A:46:VAL:HG11	2.19	0.42
1:C:164:HIS:HA	1:C:165:PRO:HD2	1.83	0.42
1:D:124:ASP:HA	1:D:149:TYR:OH	2.19	0.42
1:D:162:PRO:HG2	1:D:202:GLU:OE2	2.19	0.42
1:D:196:PHE:HE1	1:D:222:ARG:HB2	1.79	0.42
1:A:28:SER:O	1:A:29:VAL:C	2.63	0.41
1:B:14:GLU:HB2	1:B:22:MET:HE3	2.00	0.41
1:B:91:ARG:HA	1:B:129:LEU:HD21	2.01	0.41
1:C:170:LEU:O	1:C:174:PHE:HB2	2.20	0.41
1:D:78:MET:HE2	1:D:78:MET:HB3	1.96	0.41
1:B:172:LEU:HD23	2:Q:369:PHE:N	2.36	0.41
1:C:112:ALA:O	1:C:116:VAL:HG23	2.21	0.41
1:D:196:PHE:C	1:D:196:PHE:CD2	2.99	0.41
1:B:177:PHE:CE1	1:B:181:ILE:HG21	2.56	0.41
1:D:38:ASN:N	1:D:38:ASN:ND2	2.68	0.41
1:A:59:TRP:CE2	1:A:132:VAL:HG12	2.56	0.41
1:C:122:LYS:O	1:C:126:TYR:HD2	2.04	0.41
1:A:189:CYS:SG	1:A:228:TRP:HE3	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:TYR:HB3	1:B:174:PHE:CD1	2.56	0.41
1:B:167:ARG:CG	3:B:1229:PGE:H42	2.44	0.41
1:B:179:TYR:HD1	1:B:180:GLU:HG2	1.86	0.40
1:A:98:LEU:HD13	1:A:126:TYR:CE2	2.56	0.40
1:D:34:ALA:O	1:D:36:LEU:HD23	2.21	0.40
1:D:196:PHE:CD1	1:D:222:ARG:HB2	2.56	0.40
1:A:11:LYS:HE3	1:A:43:LEU:HD21	2.04	0.40
1:A:178:TYR:O	1:A:183:ASN:N	2.54	0.40
1:D:94:CYS:HB2	1:D:129:LEU:HD21	2.04	0.40
1:D:182:LEU:O	1:D:183:ASN:C	2.64	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	215/245 (88%)	195 (91%)	17 (8%)	3 (1%)	9	5
1	B	225/245 (92%)	215 (96%)	7 (3%)	3 (1%)	9	6
1	C	224/245 (91%)	211 (94%)	10 (4%)	3 (1%)	9	6
1	D	216/245 (88%)	202 (94%)	13 (6%)	1 (0%)	24	24
2	Q	7/31 (23%)	6 (86%)	1 (14%)	0	100	100
2	R	3/31 (10%)	3 (100%)	0	0	100	100
All	All	890/1042 (85%)	832 (94%)	48 (5%)	10 (1%)	11	8

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	34	ALA
1	A	109	ALA

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Mol	Chain	Res	Type
1	A	211	TYR
1	D	73	GLU
1	A	29	VAL
1	B	3	LYS
1	B	136	ASP
1	B	135	GLY
1	C	32	GLN
1	C	73	GLU

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	148/209 (71%)	145 (98%)	3 (2%)	48	59
1	B	171/209 (82%)	160 (94%)	11 (6%)	16	15
1	C	156/209 (75%)	148 (95%)	8 (5%)	21	23
1	D	140/209 (67%)	134 (96%)	6 (4%)	26	31
2	Q	3/27 (11%)	2 (67%)	1 (33%)	0	0
2	R	1/27 (4%)	1 (100%)	0	100	100
All	All	619/890 (70%)	590 (95%)	29 (5%)	23	26

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	132	VAL
1	A	170	LEU
1	A	183	ASN
1	B	52	VAL
1	B	66	GLU
1	B	78	MET
1	B	83	ARG
1	B	101	LEU
1	B	154	GLU
1	B	163	THR

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Mol	Chain	Res	Type
1	B	182	LEU
1	B	194	THR
1	B	208	GLU
1	B	216	LEU
1	C	63	SER
1	C	74	LYS
1	C	83	ARG
1	C	114	SER
1	C	154	GLU
1	C	161	GLN
1	C	182	LEU
1	C	222	ARG
1	D	51	VAL
1	D	52	VAL
1	D	203	LEU
1	D	204	ASP
1	D	219	GLN
1	D	227	LEU
2	Q	344	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	B	38	ASN
1	B	161	GLN
1	B	183	ASN
1	C	38	ASN
1	C	42	ASN
1	C	77	GLN
1	C	146	GLN
1	C	150	GLN
1	D	38	ASN
1	D	147	GLN
1	D	219	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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
2	SEP	Q	368	2	8,9,10	1.59	1 (12%)	7,12,14	1.46	2 (28%)
2	SEP	R	346	2	8,9,10	1.58	1 (12%)	7,12,14	1.47	1 (14%)
2	SEP	Q	346	2	8,9,10	1.25	1 (12%)	7,12,14	2.65	2 (28%)
2	SEP	R	368	2	8,9,10	1.63	1 (12%)	7,12,14	0.91	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	SEP	Q	368	2	-	0/6/8/10	-
2	SEP	R	346	2	-	0/6/8/10	-
2	SEP	Q	346	2	-	0/6/8/10	-
2	SEP	R	368	2	-	5/6/8/10	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	R	368	SEP	P-O1P	3.54	1.61	1.50
2	R	346	SEP	P-O1P	3.46	1.61	1.50
2	Q	368	SEP	P-O1P	3.44	1.61	1.50
2	Q	346	SEP	P-OG	-2.13	1.53	1.60

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	346	SEP	OG-CB-CA	5.84	113.83	108.14
2	R	346	SEP	OG-CB-CA	2.89	110.96	108.14
2	Q	368	SEP	OG-CB-CA	2.56	110.64	108.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	346	SEP	O3P-P-O1P	-2.51	101.05	110.83
2	Q	368	SEP	O3P-P-OG	2.41	112.95	106.67

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	R	368	SEP	N-CA-CB-OG
2	R	368	SEP	CB-OG-P-O2P
2	R	368	SEP	CB-OG-P-O3P
2	R	368	SEP	CB-OG-P-O1P
2	R	368	SEP	C-CA-CB-OG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	PGE	B	1229	-	6,6,9	0.49	0	5,5,8	0.20	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
3	PGE	B	1229	-	-	1/4/4/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1229	PGE	O2-C3-C4-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1229	PGE	2	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](#)

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	221/245 (90%)	-1.23	0 100 100	24, 34, 44, 50	0
1	B	227/245 (92%)	-1.29	0 100 100	23, 31, 40, 48	0
1	C	226/245 (92%)	-1.20	0 100 100	23, 35, 46, 56	0
1	D	222/245 (90%)	-1.09	0 100 100	29, 39, 49, 59	0
2	Q	11/31 (35%)	-1.01	0 100 100	33, 38, 43, 44	0
2	R	7/31 (22%)	-0.86	0 100 100	33, 41, 43, 43	0
All	All	914/1042 (87%)	-1.20	0 100 100	23, 35, 46, 59	0

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
2	SEP	Q	346	10/11	0.99	0.03	27,32,36,38	0
2	SEP	Q	368	10/11	0.99	0.03	21,28,32,32	0
2	SEP	R	368	10/11	0.99	0.04	31,35,40,43	0
2	SEP	R	346	10/11	1.00	0.02	20,27,32,32	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
3	PGE	B	1229	7/10	0.99	0.06	26,30,33,36	0
4	CA	D	1230	1/1	0.99	0.04	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.