



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:45 PM GMT

PDB ID : 1QJB
Title : 14-3-3 ZETA/PHOSHOPEPTIDE COMPLEX (MODE 1)
Authors : Rittinger, K.; Budman, J.; Xu, J.; Volinia, S.; Cantley, L.C.; Smerdon, S.J.;
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Deposited on : 1999-06-23
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

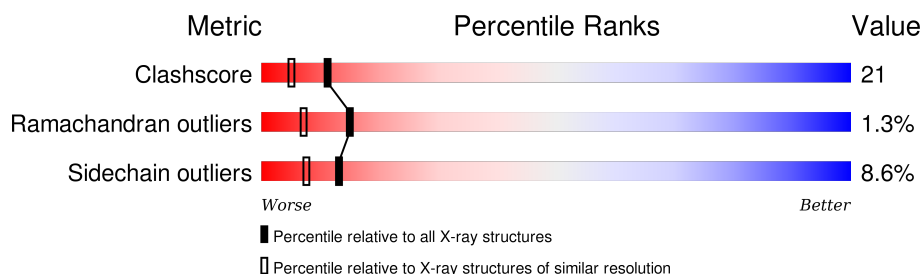
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	245	
1	B	245	
2	Q	8	
2	S	8	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4193 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	228	Total	C	N	O	S	0	0	0
			1826	1146	307	363	10			
1	B	232	Total	C	N	O	S	0	0	0
			1847	1158	310	369	10			

- Molecule 2 is a protein called PHOSPHOPEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Q	8	Total	C	N	O	P	0	0	0
			66	38	13	14	1			
2	S	7	Total	C	N	O	P	0	0	0
			61	35	12	13	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	190	Total	O	0	0
			190	190		
3	B	184	Total	O	0	0
			184	184		
3	Q	9	Total	O	0	0
			9	9		
3	S	10	Total	O	0	0
			10	10		

ALA
R4
S5
H6
S7
Y8
P9
AI0

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.35Å 71.98Å 131.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.00	Depositor
% Data completeness (in resolution range)	83.5 (15.00-2.00)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.210 , 0.288	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4193	wwPDB-VP
Average B, all atoms (Å ²)	23.0	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: SEP

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	0.59	0/1850	1.45	25/2486 (1.0%)
1	B	0.61	0/1872	1.38	18/2518 (0.7%)
2	Q	0.66	0/57	1.55	0/75
2	S	0.77	0/52	1.95	3/68 (4.4%)
All	All	0.61	0/3831	1.43	46/5147 (0.9%)

There are no bond length outliers.

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	ARG	NE-CZ-NH1	13.48	127.04	120.30
1	B	60	ARG	NE-CZ-NH1	13.14	126.87	120.30
1	A	41	ARG	NE-CZ-NH1	10.17	125.39	120.30
1	B	83	ARG	NE-CZ-NH1	9.89	125.25	120.30
1	A	60	ARG	NE-CZ-NH1	9.86	125.23	120.30
1	A	222	ARG	CD-NE-CZ	9.08	136.32	123.60
1	A	60	ARG	NE-CZ-NH2	-9.05	115.77	120.30
1	B	83	ARG	CD-NE-CZ	8.74	135.84	123.60
1	A	83	ARG	CD-NE-CZ	8.71	135.79	123.60
1	B	127	ARG	NE-CZ-NH2	-8.37	116.12	120.30
1	B	60	ARG	CD-NE-CZ	8.33	135.26	123.60
1	A	140	GLY	C-N-CA	8.10	141.95	121.70
1	A	127	ARG	NE-CZ-NH1	7.99	124.30	120.30
1	A	80	ARG	CD-NE-CZ	7.41	133.97	123.60
1	B	83	ARG	NE-CZ-NH2	-7.39	116.60	120.30
1	A	83	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	B	140	GLY	O-C-N	-7.11	111.32	122.70
1	B	60	ARG	NE-CZ-NH2	-6.96	116.82	120.30
1	A	167	ARG	NE-CZ-NH2	-6.82	116.89	120.30
1	A	124	ASP	CB-CG-OD1	6.78	124.40	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	141	ILE	N-CA-CB	6.69	126.19	110.80
2	S	8	TYR	CA-C-O	-6.65	106.13	120.10
1	B	20	ASP	CB-CG-OD1	6.57	124.21	118.30
1	A	215	THR	N-CA-CB	-6.42	98.11	110.30
1	A	18	ARG	NE-CZ-NH2	-6.32	117.14	120.30
1	B	104	PHE	CB-CA-C	-6.27	97.86	110.40
1	B	223	ASP	CB-CG-OD1	6.25	123.93	118.30
2	S	9	PRO	N-CA-CB	6.13	110.65	103.30
1	A	141	ILE	CA-C-O	5.96	132.63	120.10
1	A	132	VAL	N-CA-CB	-5.96	98.40	111.50
1	B	222	ARG	NE-CZ-NH2	-5.85	117.38	120.30
1	A	18	ARG	NE-CZ-NH1	5.78	123.19	120.30
2	S	8	TYR	O-C-N	5.76	132.04	121.10
1	A	55	ARG	NE-CZ-NH2	-5.75	117.42	120.30
1	B	1	MET	CG-SD-CE	5.67	109.27	100.20
1	B	55	ARG	NE-CZ-NH2	-5.61	117.50	120.30
1	A	41	ARG	CD-NE-CZ	5.58	131.41	123.60
1	A	91	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	A	83	ARG	CG-CD-NE	5.37	123.08	111.80
1	B	109	ALA	CB-CA-C	5.32	118.07	110.10
1	A	118	TYR	CB-CG-CD2	-5.27	117.84	121.00
1	A	203	LEU	C-N-CA	5.20	134.70	121.70
1	B	140	GLY	N-CA-C	5.18	126.05	113.10
1	B	149	TYR	CB-CG-CD2	-5.14	117.92	121.00
1	A	225	LEU	CA-CB-CG	5.11	127.04	115.30
1	B	139	LYS	CA-C-N	5.02	126.24	116.20

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	1826	0	1812	87	0
1	B	1847	0	1828	66	1
2	Q	66	0	53	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	S	61	0	48	3	0
3	A	190	0	0	55	0
3	B	184	0	0	35	1
3	Q	9	0	0	3	0
3	S	10	0	0	3	0
All	All	4193	0	3741	156	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 21.

All (156) 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:140:GLY:HA2	3:B:2126:HOH:O	1.09	1.23
1:A:140:GLY:HA2	3:A:2128:HOH:O	1.06	1.21
1:B:42:ASN:HB3	3:B:2042:HOH:O	1.38	1.19
1:A:144:GLN:HG3	3:A:2135:HOH:O	1.41	1.16
1:A:38:ASN:HB2	3:A:2047:HOH:O	1.44	1.12
1:B:139:LYS:CE	3:B:2122:HOH:O	1.98	1.09
1:B:209:GLU:HG3	3:B:2171:HOH:O	1.53	1.08
1:B:80:ARG:HD2	3:B:2081:HOH:O	1.51	1.06
1:A:91:ARG:NH1	3:A:2096:HOH:O	1.92	1.00
1:B:130:ALA:HA	1:B:141:ILE:HD11	1.44	0.98
1:A:139:LYS:HE3	3:A:2127:HOH:O	1.61	0.98
1:A:191:LEU:HD13	3:A:2033:HOH:O	1.67	0.94
1:B:139:LYS:CG	3:B:2122:HOH:O	2.15	0.93
1:A:101:LEU:HD21	1:A:121:MET:HE2	1.51	0.93
1:A:141:ILE:HA	3:A:2129:HOH:O	1.69	0.92
1:A:140:GLY:CA	3:A:2128:HOH:O	1.76	0.92
1:B:4:ASN:O	3:B:2004:HOH:O	1.88	0.92
1:A:139:LYS:CE	3:A:2127:HOH:O	2.18	0.91
1:A:78:MET:HB3	3:A:2086:HOH:O	1.68	0.90
1:B:139:LYS:HE3	3:B:2122:HOH:O	1.67	0.87
1:A:38:ASN:CB	3:A:2047:HOH:O	2.08	0.86
1:A:41:ARG:NE	3:A:2053:HOH:O	2.09	0.86
1:A:8:GLN:HG2	3:A:2005:HOH:O	1.76	0.84
1:A:41:ARG:CD	3:A:2053:HOH:O	2.24	0.84
1:A:73:GLU:OE2	3:A:2081:HOH:O	1.93	0.84
1:A:139:LYS:HD3	3:A:2131:HOH:O	1.79	0.83
1:B:164:HIS:NE2	3:B:2142:HOH:O	2.12	0.82
1:A:8:GLN:HB2	3:A:2007:HOH:O	1.79	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:ILE:CA	3:A:2129:HOH:O	2.25	0.81
1:B:85:LYS:NZ	3:B:2084:HOH:O	2.13	0.81
1:A:63:SER:O	1:A:67:GLN:HG3	1.82	0.79
1:A:4:ASN:HB3	3:A:2003:HOH:O	1.84	0.78
1:B:139:LYS:HG3	3:B:2122:HOH:O	1.75	0.77
1:A:216:LEU:HD12	3:A:2181:HOH:O	1.86	0.76
1:A:78:MET:N	3:A:2086:HOH:O	2.20	0.75
1:A:180:GLU:HG3	3:A:2158:HOH:O	1.86	0.75
1:A:38:ASN:ND2	3:A:2047:HOH:O	2.18	0.75
1:B:38:ASN:CB	3:B:2039:HOH:O	2.36	0.74
1:B:191:LEU:HD12	3:B:2159:HOH:O	1.87	0.74
1:B:138:LYS:HE3	3:B:2121:HOH:O	1.87	0.74
1:A:140:GLY:C	3:A:2128:HOH:O	2.05	0.73
1:A:191:LEU:HD12	3:A:2167:HOH:O	1.89	0.72
1:A:38:ASN:HD22	1:A:41:ARG:HH22	1.36	0.72
1:B:113:GLU:HA	1:B:160:MET:HE2	1.70	0.72
1:B:139:LYS:NZ	3:B:2122:HOH:O	1.99	0.71
1:A:130:ALA:HA	1:A:141:ILE:HD11	1.73	0.71
1:A:38:ASN:HD22	1:A:41:ARG:NH2	1.89	0.71
1:B:38:ASN:HB2	3:B:2039:HOH:O	1.90	0.70
1:B:130:ALA:CA	1:B:141:ILE:HD11	2.20	0.70
1:B:38:ASN:ND2	3:B:2039:HOH:O	1.94	0.70
1:B:60:ARG:NE	3:B:2063:HOH:O	2.24	0.69
1:B:136:ASP:O	1:B:137:ASP:HB2	1.94	0.67
1:A:139:LYS:HB2	3:A:2123:HOH:O	1.94	0.67
1:B:2:ASP:OD1	1:B:3:LYS:N	2.28	0.67
1:A:8:GLN:CB	3:A:2007:HOH:O	2.37	0.67
1:A:211:TYR:O	1:A:215:THR:HB	1.94	0.67
1:B:140:GLY:CA	3:B:2126:HOH:O	1.91	0.66
1:B:137:ASP:O	3:B:2121:HOH:O	2.15	0.65
1:B:164:HIS:CD2	3:B:2142:HOH:O	2.49	0.65
1:B:225:LEU:O	1:B:229:THR:HB	1.97	0.64
1:A:139:LYS:CB	3:A:2123:HOH:O	2.44	0.64
1:A:138:LYS:HE2	1:A:181:ILE:HA	1.79	0.64
1:B:113:GLU:HA	1:B:160:MET:CE	2.29	0.62
1:A:139:LYS:HG3	1:A:140:GLY:N	2.13	0.62
1:A:135:GLY:O	1:A:136:ASP:HB2	2.00	0.62
1:B:191:LEU:HD13	3:B:2026:HOH:O	1.99	0.62
1:A:157:LYS:CE	3:A:2147:HOH:O	2.47	0.62
1:A:41:ARG:NH2	3:A:2052:HOH:O	2.30	0.61
1:B:102:GLU:HG2	3:B:2097:HOH:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:GLU:HA	1:A:160:MET:CE	2.31	0.61
1:B:164:HIS:CE1	3:B:2142:HOH:O	2.52	0.61
2:Q:4:ARG:HG2	3:Q:2003:HOH:O	2.02	0.60
1:B:4:ASN:O	1:B:8:GLN:HG3	2.02	0.59
1:B:60:ARG:CZ	3:B:2063:HOH:O	2.49	0.59
1:A:135:GLY:HA2	3:A:2124:HOH:O	2.02	0.59
2:Q:4:ARG:CG	3:Q:2003:HOH:O	2.50	0.59
2:S:6:HIS:CE1	3:S:2003:HOH:O	2.55	0.58
1:B:196:PHE:CZ	1:B:200:ILE:HD12	2.39	0.57
1:A:106:ILE:N	1:A:107:PRO:HD2	2.20	0.57
1:A:8:GLN:HG3	3:A:2007:HOH:O	2.05	0.56
1:B:139:LYS:HD3	1:B:140:GLY:N	2.20	0.56
1:A:41:ARG:NE	3:A:2052:HOH:O	2.06	0.56
1:B:59:TRP:CE2	1:B:132:VAL:HG12	2.41	0.56
1:B:9:LYS:HD3	1:B:25:CYS:SG	2.46	0.56
1:B:138:LYS:HE2	1:B:181:ILE:O	2.05	0.55
1:B:33:GLY:HA2	1:B:104:PHE:CE1	2.42	0.55
1:A:180:GLU:CG	3:A:2158:HOH:O	2.51	0.54
1:A:153:PHE:CZ	1:A:191:LEU:HD11	2.43	0.54
1:A:157:LYS:HE2	3:A:2147:HOH:O	2.05	0.54
1:A:139:LYS:HE3	3:A:2123:HOH:O	2.07	0.54
1:A:41:ARG:NH2	3:A:2047:HOH:O	2.40	0.54
1:B:4:ASN:HB2	3:B:2001:HOH:O	2.07	0.54
1:A:111:GLN:HB2	3:A:2104:HOH:O	2.07	0.53
1:B:139:LYS:O	1:B:141:ILE:N	2.41	0.53
1:A:59:TRP:CE2	1:A:132:VAL:HG22	2.43	0.53
1:B:139:LYS:HD2	1:B:141:ILE:HG23	1.91	0.53
1:A:78:MET:CA	3:A:2086:HOH:O	2.57	0.53
1:A:87:GLU:OE2	1:A:91:ARG:NH2	2.42	0.52
1:B:116:VAL:HG21	1:B:160:MET:HE3	1.91	0.52
1:A:41:ARG:HD3	3:A:2053:HOH:O	1.99	0.52
1:B:89:GLU:OE2	3:B:2087:HOH:O	2.19	0.52
1:A:143:ASP:HB3	3:A:2133:HOH:O	2.10	0.52
1:A:78:MET:CB	3:A:2086:HOH:O	2.39	0.52
1:A:225:LEU:O	1:A:229:THR:HB	2.09	0.51
1:A:130:ALA:CA	1:A:141:ILE:HD11	2.40	0.51
1:A:231:ASP:C	3:A:2190:HOH:O	2.50	0.50
1:A:231:ASP:CA	3:A:2190:HOH:O	2.59	0.50
1:B:109:ALA:O	1:B:110:SER:CB	2.60	0.50
1:A:153:PHE:CZ	1:A:157:LYS:HE3	2.47	0.49
2:S:4:ARG:N	3:S:2002:HOH:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:GLN:HB3	1:B:78:MET:SD	2.53	0.49
1:B:94:CYS:HB2	1:B:129:LEU:HD21	1.95	0.49
1:A:113:GLU:HA	1:A:160:MET:HE2	1.95	0.48
1:A:139:LYS:HB3	3:A:2123:HOH:O	2.11	0.48
1:A:139:LYS:O	1:A:141:ILE:HD13	2.14	0.48
1:A:12:LEU:HD12	1:B:78:MET:HG2	1.96	0.47
1:A:216:LEU:CD1	3:A:2181:HOH:O	2.52	0.47
1:A:80:ARG:HD3	3:A:2087:HOH:O	2.14	0.47
1:B:206:LEU:HD13	1:B:211:TYR:CE1	2.48	0.47
1:B:80:ARG:O	1:B:84:GLU:HG3	2.15	0.47
1:B:113:GLU:HG2	3:B:2142:HOH:O	2.13	0.46
1:A:8:GLN:NE2	3:A:2006:HOH:O	2.46	0.46
1:A:121:MET:HG2	3:A:2058:HOH:O	2.15	0.46
1:B:110:SER:OG	1:B:111:GLN:NE2	2.48	0.46
1:A:139:LYS:HD2	3:A:2132:HOH:O	2.16	0.46
1:B:160:MET:HG2	3:B:2142:HOH:O	2.15	0.45
1:B:49:LYS:HD3	3:B:2112:HOH:O	2.16	0.45
1:A:8:GLN:CG	3:A:2007:HOH:O	2.60	0.45
1:A:38:ASN:ND2	1:A:41:ARG:HH22	2.11	0.45
1:B:143:ASP:O	1:B:147:GLN:HG3	2.16	0.44
1:A:141:ILE:HD13	1:A:142:VAL:N	2.33	0.44
1:A:139:LYS:HG2	1:A:141:ILE:HG23	2.00	0.44
1:A:222:ARG:HG3	1:A:222:ARG:HH11	1.83	0.44
1:A:1:MET:CE	1:A:9:LYS:HD2	2.48	0.44
1:B:8:GLN:HG3	3:B:2004:HOH:O	2.17	0.44
1:A:177:PHE:CE2	1:A:182:LEU:HD22	2.53	0.43
1:A:30:THR:HG21	1:A:100:LEU:HG	2.00	0.43
2:Q:4:ARG:HG3	3:Q:2003:HOH:O	2.16	0.43
1:B:211:TYR:O	1:B:215:THR:HB	2.18	0.43
1:B:138:LYS:O	1:B:139:LYS:HB2	2.18	0.42
1:B:139:LYS:O	1:B:141:ILE:HD13	2.19	0.42
1:A:164:HIS:HE1	1:A:166:ILE:HD12	1.83	0.42
1:A:167:ARG:NH2	3:A:2153:HOH:O	2.46	0.42
1:B:26:MET:O	1:B:30:THR:HG23	2.20	0.42
1:A:135:GLY:CA	3:A:2124:HOH:O	2.65	0.42
1:A:141:ILE:HD13	1:A:142:VAL:H	1.85	0.41
1:A:37:SER:OG	1:A:40:GLU:HG3	2.20	0.41
2:S:10:ALA:C	3:S:2010:HOH:O	2.59	0.41
1:B:56:ARG:HH11	1:B:56:ARG:HD3	1.68	0.41
1:B:139:LYS:HG2	3:B:2122:HOH:O	2.03	0.41
1:A:183:ASN:ND2	3:A:2159:HOH:O	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:ASP:CG	3:B:2001:HOH:O	2.59	0.40
1:B:140:GLY:C	3:B:2126:HOH:O	2.44	0.40
1:A:76:GLN:C	3:A:2086:HOH:O	2.59	0.40
1:A:12:LEU:CD1	1:B:78:MET:HG2	2.52	0.40
1:B:210:SER:HA	3:B:2176:HOH:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:LYS:CG	3:B:2170:HOH:O[4_456]	2.18	0.02

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	224/245 (91%)	215 (96%)	7 (3%)	2 (1%)	21	13
1	B	230/245 (94%)	217 (94%)	9 (4%)	4 (2%)	11	4
2	Q	5/8 (62%)	5 (100%)	0	0	100	100
2	S	4/8 (50%)	4 (100%)	0	0	100	100
All	All	463/506 (92%)	441 (95%)	16 (4%)	6 (1%)	15	7

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	B	137	ASP
1	B	140	GLY
1	A	140	GLY
1	B	231	ASP

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Mol	Chain	Res	Type
1	B	109	ALA

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	197/209 (94%)	180 (91%)	17 (9%)	13	7
1	B	198/209 (95%)	181 (91%)	17 (9%)	13	7
2	Q	5/5 (100%)	5 (100%)	0	100	100
2	S	5/5 (100%)	4 (80%)	1 (20%)	1	0
All	All	405/428 (95%)	370 (91%)	35 (9%)	13	7

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	8	GLN
1	A	63	SER
1	A	66	GLU
1	A	100	LEU
1	A	108	ASN
1	A	115	LYS
1	A	136	ASP
1	A	139	LYS
1	A	141	ILE
1	A	174	PHE
1	A	182	LEU
1	A	183	ASN
1	A	211	TYR
1	A	215	THR
1	A	222	ARG
1	A	225	LEU
1	B	1	MET
1	B	3	LYS
1	B	30	THR

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Mol	Chain	Res	Type
1	B	41	ARG
1	B	100	LEU
1	B	103	LYS
1	B	110	SER
1	B	114	SER
1	B	136	ASP
1	B	139	LYS
1	B	141	ILE
1	B	157	LYS
1	B	174	PHE
1	B	200	ILE
1	B	204	ASP
1	B	215	THR
1	B	229	THR
2	S	4	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	76	GLN
1	B	108	ASN
1	B	111	GLN
1	B	183	ASN
1	B	219	GLN

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
2	SEP	Q	7	2	8,9,10	3.50	1 (12%)	8,12,14	6.18	4 (50%)
2	SEP	S	7	2	8,9,10	3.65	1 (12%)	8,12,14	3.80	2 (25%)

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	7	2	-	0/6/8/10	0/0/0/0
2	SEP	S	7	2	-	0/6/8/10	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	7	SEP	P-OG	9.71	1.92	1.60
2	S	7	SEP	P-OG	10.12	1.94	1.60

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	7	SEP	OG-CB-CA	-16.46	94.23	108.27
2	S	7	SEP	OG-CB-CA	-9.62	100.07	108.27
2	Q	7	SEP	O-C-CA	-2.85	118.08	125.49
2	Q	7	SEP	O2P-P-O1P	-2.08	103.88	110.58
2	S	7	SEP	O3P-P-OG	3.54	116.76	106.56
2	Q	7	SEP	O3P-P-OG	4.17	118.58	106.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.