



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:30 AM UTC

PDB ID : 1A37 / pdb_00001a37
Title : 14-3-3 PROTEIN ZETA BOUND TO PS-RAF259 PEPTIDE
Authors : Petosa, C.; Masters, S.C.; Pohl, J.; Wang, B.; Fu, H.; Liddington, R.C.
Deposited on : 1998-01-28
Resolution : 3.60 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

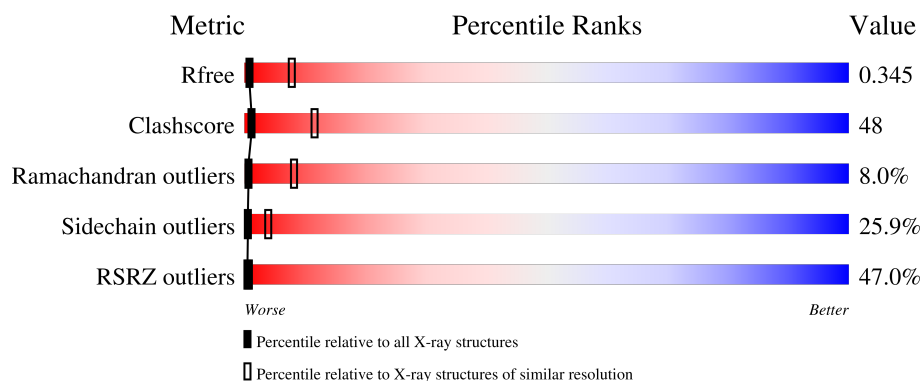
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	A	245	
1	B	245	
2	P	15	
2	Q	15	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3194 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1563	984	265	305	9			
1	B	195	Total	C	N	O	S	0	0	0
			1563	984	265	305	9			

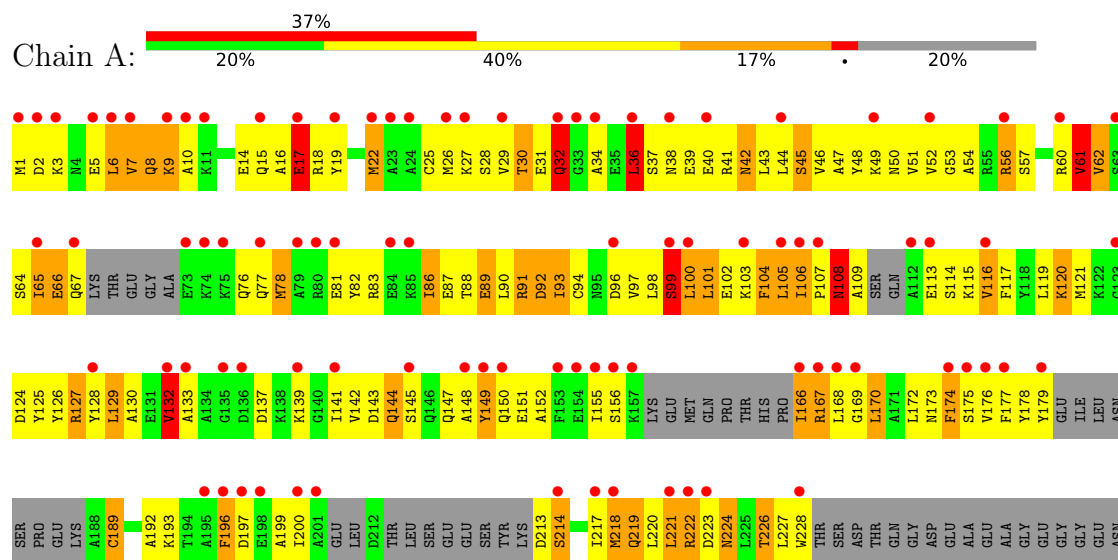
- Molecule 2 is a protein called PS-RAF259 PEPTIDE LSQRQRST(SEP)TPNVHM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	7	Total	C	N	O	P	4	0	0
			34	15	7	11	1			
2	Q	7	Total	C	N	O	P	4	0	0
			34	15	7	11	1			

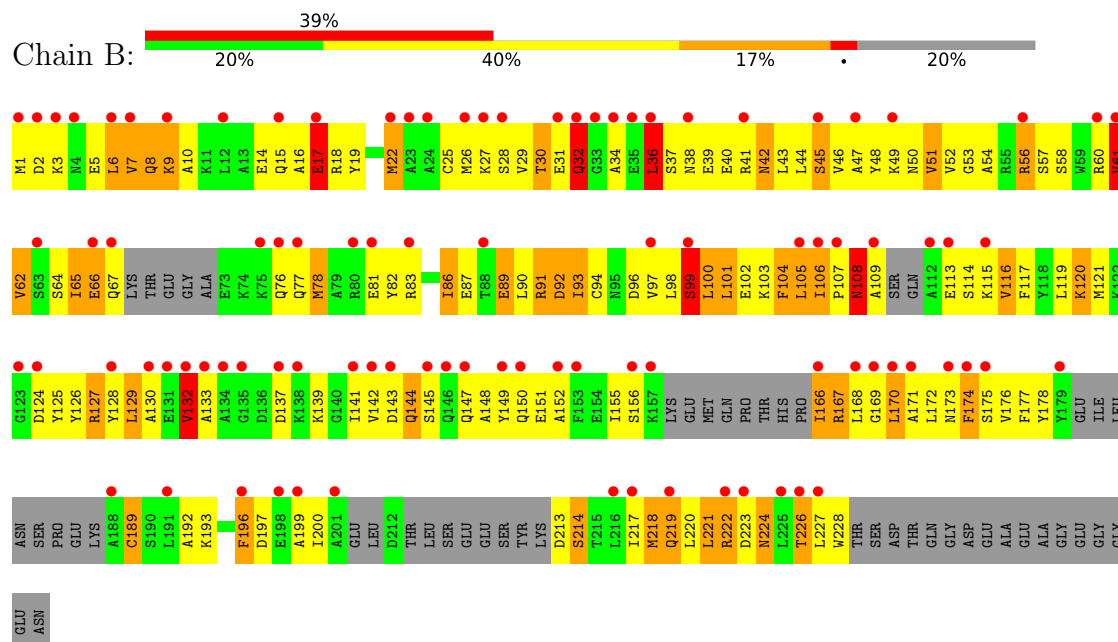
3 Residue-property plots

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

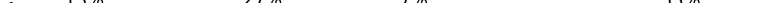
• Molecule 1: 14-3-3 PROTEIN ZETA



• Molecule 1: 14-3-3 PROTEIN ZETA

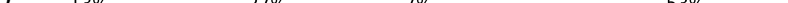


• Molecule 2: PS-RAF259 PEPTIDE LSQRQRST(SEP)TPNVHM

Chain P:  13% 27% 7% 53%



- Molecule 2: PS-RAF259 PEPTIDE LSQRQRST(SEP)TPNVHM

Chain Q: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	94.73Å 94.73Å 250.87Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.60 20.00 – 3.60	Depositor EDS
% Data completeness (in resolution range)	98.9 (20.00-3.60) 96.5 (20.00-3.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.47 (at 3.32Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.320 , 0.360 0.322 , 0.345	Depositor DCC
R_{free} test set	1873 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å ²)	76.5	Xtriage
Anisotropy	0.904	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.115 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.64	EDS
Total number of atoms	3194	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.56% 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

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	0.82	2/1578 (0.1%)	1.35	15/2111 (0.7%)
1	B	0.83	2/1578 (0.1%)	1.35	15/2111 (0.7%)
2	P	0.90	0/22	1.59	0/24
2	Q	0.91	0/22	1.59	0/24
All	All	0.83	4/3200 (0.1%)	1.35	30/4270 (0.7%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	106	ILE	CA-CB	13.00	1.60	1.54
1	A	106	ILE	CA-CB	12.87	1.60	1.54
1	A	218	MET	SD-CE	5.08	1.92	1.79
1	B	218	MET	SD-CE	5.08	1.92	1.79

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	132	VAL	N-CA-C	9.14	119.95	110.72
1	B	132	VAL	N-CA-C	9.14	119.95	110.72
1	B	127	ARG	N-CA-C	-8.32	102.16	111.07
1	A	127	ARG	N-CA-C	-8.30	102.19	111.07
1	B	167	ARG	N-CA-C	-7.62	102.11	111.33
1	A	167	ARG	N-CA-C	-7.60	102.14	111.33
1	A	32	GLN	N-CA-C	-7.58	94.66	110.80
1	A	226	THR	N-CA-C	-7.58	102.16	111.33
1	B	32	GLN	N-CA-C	-7.58	94.66	110.80
1	B	226	THR	N-CA-C	-7.56	102.18	111.33
1	B	7	VAL	N-CA-C	-7.19	104.20	110.74
1	A	7	VAL	N-CA-C	-7.17	104.22	110.74
1	A	77	GLN	N-CA-C	-6.79	103.91	111.71

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	77	GLN	N-CA-C	-6.78	103.91	111.71
1	A	222	ARG	N-CA-C	-6.50	103.89	110.97
1	B	222	ARG	N-CA-C	-6.50	103.89	110.97
1	A	6	LEU	N-CA-C	-6.39	104.97	112.89
1	B	6	LEU	N-CA-C	-6.38	104.98	112.89
1	B	36	LEU	N-CA-C	6.19	119.34	109.24
1	A	36	LEU	N-CA-C	6.18	119.32	109.24
1	B	174	PHE	N-CA-C	-6.02	104.63	111.07
1	A	174	PHE	N-CA-C	-6.01	104.64	111.07
1	B	61	VAL	N-CA-C	5.94	116.41	110.23
1	A	61	VAL	N-CA-C	5.92	116.39	110.23
1	A	120	LYS	N-CA-C	-5.49	104.94	111.03
1	B	120	LYS	N-CA-C	-5.48	104.95	111.03
1	B	17	GLU	N-CA-C	5.35	122.20	110.80
1	A	17	GLU	N-CA-C	5.34	122.17	110.80
1	A	106	ILE	CB-CA-C	5.16	118.60	113.70
1	B	106	ILE	CB-CA-C	5.13	118.57	113.70

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	1563	0	1558	155	0
1	B	1563	0	1558	153	0
2	P	34	0	7	1	0
2	Q	34	0	7	1	0
All	All	3194	0	3130	304	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 48.

All (304) 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:A:22:MET:SD	1:A:47:ALA:HB2	2.09	0.93
1:B:22:MET:SD	1:B:47:ALA:HB2	2.09	0.93
1:A:46:VAL:HA	1:A:49:LYS:HG2	1.63	0.79
1:B:46:VAL:HA	1:B:49:LYS:HG2	1.63	0.78
1:A:15:GLN:O	1:B:61:VAL:HG21	1.83	0.78
1:A:94:CYS:SG	1:A:129:LEU:HD11	2.24	0.78
1:B:94:CYS:SG	1:B:129:LEU:HD11	2.24	0.77
1:A:92:ASP:O	1:A:96:ASP:HB2	1.84	0.77
1:B:92:ASP:O	1:B:96:ASP:HB2	1.85	0.77
1:B:94:CYS:SG	1:B:129:LEU:CD1	2.73	0.77
1:B:101:LEU:HA	1:B:105:LEU:HB2	1.66	0.76
1:A:94:CYS:SG	1:A:129:LEU:CD1	2.73	0.76
1:A:101:LEU:HA	1:A:105:LEU:HB2	1.66	0.76
1:A:48:TYR:O	1:A:52:VAL:HG12	1.86	0.75
1:B:48:TYR:O	1:B:52:VAL:HG12	1.86	0.74
1:B:172:LEU:O	1:B:175:SER:HB3	1.87	0.73
1:A:223:ASP:O	1:A:226:THR:HB	1.89	0.73
1:A:172:LEU:O	1:A:175:SER:HB3	1.87	0.73
1:B:223:ASP:O	1:B:226:THR:HB	1.89	0.73
1:B:30:THR:HG21	1:B:100:LEU:HD23	1.71	0.72
1:A:87:GLU:HG3	1:A:132:VAL:CG1	2.20	0.71
1:A:30:THR:HG21	1:A:100:LEU:HD23	1.71	0.71
1:B:87:GLU:HG3	1:B:132:VAL:CG1	2.20	0.70
1:A:61:VAL:HG21	1:B:15:GLN:O	1.91	0.70
1:B:214:SER:O	1:B:218:MET:HG3	1.92	0.70
1:A:89:GLU:O	1:A:93:ILE:HG12	1.92	0.70
1:A:214:SER:O	1:A:218:MET:HG3	1.92	0.69
1:B:89:GLU:O	1:B:93:ILE:HG12	1.92	0.69
1:B:223:ASP:O	1:B:227:LEU:HG	1.93	0.68
1:A:221:LEU:H	1:A:221:LEU:HD23	1.59	0.68
1:B:221:LEU:HD23	1:B:221:LEU:H	1.59	0.68
1:A:124:ASP:O	1:A:127:ARG:HB3	1.94	0.67
1:A:223:ASP:O	1:A:227:LEU:HG	1.93	0.67
1:B:38:ASN:OD1	1:B:41:ARG:HD3	1.94	0.67
1:B:124:ASP:O	1:B:127:ARG:HB3	1.94	0.67
1:B:166:ILE:HG23	1:B:167:ARG:H	1.60	0.66
1:A:38:ASN:OD1	1:A:41:ARG:HD3	1.94	0.66
1:B:22:MET:SD	1:B:47:ALA:CB	2.83	0.66
1:B:156:SER:HB2	1:B:167:ARG:HG3	1.77	0.66
1:A:48:TYR:CE2	1:A:97:VAL:HB	2.31	0.66
1:B:48:TYR:CE2	1:B:97:VAL:HB	2.31	0.66
1:A:93:ILE:O	1:A:97:VAL:HG12	1.97	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:GLU:HG2	1:A:91:ARG:NH2	2.12	0.65
1:A:121:MET:O	1:A:124:ASP:HB2	1.97	0.65
1:A:22:MET:SD	1:A:47:ALA:CB	2.84	0.65
1:A:166:ILE:HG23	1:A:167:ARG:H	1.60	0.65
1:B:87:GLU:HG2	1:B:91:ARG:NH2	2.12	0.65
1:B:93:ILE:O	1:B:97:VAL:HG12	1.97	0.64
1:A:156:SER:HB2	1:A:167:ARG:HG3	1.77	0.64
1:A:114:SER:O	1:A:117:PHE:HB3	1.97	0.64
1:B:114:SER:O	1:B:117:PHE:HB3	1.97	0.64
1:B:121:MET:O	1:B:124:ASP:HB2	1.97	0.64
1:A:139:LYS:O	1:A:142:VAL:HG12	1.97	0.64
1:B:217:ILE:O	1:B:221:LEU:HD23	1.98	0.64
1:A:217:ILE:O	1:A:221:LEU:HD23	1.98	0.64
1:A:31:GLU:HA	1:A:104:PHE:CZ	2.33	0.64
1:B:139:LYS:O	1:B:142:VAL:HG12	1.97	0.63
1:B:31:GLU:HA	1:B:104:PHE:CZ	2.33	0.62
1:A:217:ILE:O	1:A:220:LEU:HB2	2.00	0.62
1:B:14:GLU:C	1:B:16:ALA:H	2.08	0.62
1:A:14:GLU:C	1:A:16:ALA:H	2.08	0.62
1:A:39:GLU:O	1:A:43:LEU:HD13	2.00	0.62
1:B:39:GLU:O	1:B:43:LEU:HD13	2.00	0.62
1:B:217:ILE:O	1:B:220:LEU:HB2	2.00	0.61
1:B:98:LEU:O	1:B:102:GLU:HG3	2.01	0.61
1:B:46:VAL:HG23	1:B:49:LYS:CE	2.31	0.61
1:B:174:PHE:CD1	1:B:178:TYR:HE1	2.19	0.61
1:A:98:LEU:O	1:A:102:GLU:HG3	2.01	0.61
1:B:52:VAL:HG11	1:B:125:TYR:HE1	1.66	0.61
1:A:46:VAL:HG23	1:A:49:LYS:CE	2.31	0.60
1:A:174:PHE:CD1	1:A:178:TYR:HE1	2.19	0.60
1:A:52:VAL:HG11	1:A:125:TYR:HE1	1.66	0.60
1:A:119:LEU:HB3	1:A:152:ALA:HB2	1.82	0.60
1:B:90:LEU:HD21	1:B:132:VAL:HG21	1.83	0.60
1:B:46:VAL:HG23	1:B:49:LYS:HE2	1.84	0.60
1:A:94:CYS:SG	1:A:129:LEU:HD13	2.42	0.60
1:A:90:LEU:HD21	1:A:132:VAL:HG21	1.83	0.60
1:B:119:LEU:HB3	1:B:152:ALA:HB2	1.82	0.59
1:A:46:VAL:HG23	1:A:49:LYS:HE2	1.84	0.59
1:A:87:GLU:O	1:A:90:LEU:HB3	2.02	0.59
1:B:169:GLY:O	1:B:172:LEU:HB3	2.03	0.59
1:A:147:GLN:O	1:A:151:GLU:HB2	2.02	0.59
1:B:147:GLN:O	1:B:151:GLU:HB2	2.02	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:GLU:O	1:B:90:LEU:HB3	2.02	0.59
1:A:127:ARG:HH22	2:P:259:SEP:P	2.25	0.59
1:A:169:GLY:O	1:A:172:LEU:HB3	2.03	0.59
1:A:27:LYS:O	1:A:31:GLU:HG2	2.03	0.58
1:B:27:LYS:O	1:B:31:GLU:HG2	2.03	0.58
1:B:100:LEU:O	1:B:104:PHE:HB2	2.03	0.58
1:A:126:TYR:CE2	1:A:144:GLN:HB3	2.38	0.58
1:B:127:ARG:HH22	2:Q:259:SEP:P	2.26	0.58
1:A:100:LEU:O	1:A:104:PHE:HB2	2.03	0.58
1:B:126:TYR:CE2	1:B:144:GLN:HB3	2.38	0.58
1:B:94:CYS:SG	1:B:129:LEU:HD13	2.42	0.58
1:B:106:ILE:HG13	1:B:107:PRO:HD3	1.86	0.57
1:B:128:TYR:O	1:B:132:VAL:HG23	2.04	0.57
1:A:196:PHE:CZ	1:A:200:ILE:HD11	2.40	0.56
1:B:98:LEU:HD11	1:B:126:TYR:CE1	2.40	0.56
1:A:106:ILE:HG13	1:A:107:PRO:HD3	1.86	0.56
1:A:109:ALA:CB	1:A:115:LYS:HG3	2.36	0.56
1:A:128:TYR:O	1:A:132:VAL:HG23	2.04	0.56
1:B:150:GLN:HA	1:B:174:PHE:CE2	2.41	0.56
1:B:196:PHE:CZ	1:B:200:ILE:HD11	2.40	0.56
1:A:98:LEU:HD11	1:A:126:TYR:CE1	2.40	0.56
1:B:109:ALA:CB	1:B:115:LYS:HG3	2.36	0.56
1:A:26:MET:HB3	1:A:44:LEU:CD1	2.36	0.55
1:A:9:LYS:HG2	1:B:78:MET:SD	2.47	0.55
1:B:57:SER:O	1:B:60:ARG:HB2	2.07	0.55
1:A:193:LYS:HA	1:A:196:PHE:HD2	1.71	0.55
1:A:150:GLN:HA	1:A:174:PHE:CE2	2.41	0.55
1:A:174:PHE:O	1:A:177:PHE:HB3	2.07	0.55
1:B:26:MET:HB3	1:B:44:LEU:CD1	2.36	0.55
1:B:193:LYS:HA	1:B:196:PHE:HD2	1.71	0.55
1:B:196:PHE:HB2	1:B:221:LEU:HD12	1.89	0.55
1:A:57:SER:O	1:A:60:ARG:HB2	2.07	0.55
1:A:196:PHE:HB2	1:A:221:LEU:HD12	1.89	0.55
1:A:196:PHE:CG	1:A:197:ASP:N	2.76	0.54
1:B:174:PHE:O	1:B:177:PHE:HB3	2.07	0.54
1:B:196:PHE:CG	1:B:197:ASP:N	2.76	0.53
1:A:78:MET:SD	1:B:9:LYS:HG2	2.49	0.53
1:A:150:GLN:HA	1:A:174:PHE:HE2	1.74	0.53
1:B:127:ARG:O	1:B:130:ALA:HB3	2.09	0.53
1:B:150:GLN:HA	1:B:174:PHE:HE2	1.73	0.53
1:B:117:PHE:O	1:B:120:LYS:HB3	2.09	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:PHE:CE1	1:A:200:ILE:HD11	2.45	0.52
1:B:41:ARG:HG2	1:B:117:PHE:HE2	1.75	0.52
1:A:117:PHE:O	1:A:120:LYS:HB3	2.09	0.52
1:A:127:ARG:O	1:A:130:ALA:HB3	2.09	0.52
1:B:174:PHE:CE1	1:B:178:TYR:CE1	2.98	0.51
1:A:174:PHE:CE1	1:A:178:TYR:CE1	2.98	0.51
1:B:196:PHE:CE1	1:B:200:ILE:HD11	2.45	0.51
1:A:41:ARG:HG2	1:A:117:PHE:HE2	1.75	0.51
1:B:113:GLU:O	1:B:116:VAL:HG23	2.11	0.51
1:A:26:MET:HB3	1:A:44:LEU:HD13	1.93	0.51
1:B:109:ALA:HB1	1:B:115:LYS:HG3	1.94	0.50
1:B:224:ASN:HB3	1:B:228:TRP:CZ3	2.47	0.50
1:A:98:LEU:HD11	1:A:126:TYR:CD1	2.46	0.50
1:A:113:GLU:O	1:A:116:VAL:HG23	2.11	0.50
1:A:167:ARG:O	1:A:170:LEU:HB3	2.12	0.50
1:A:224:ASN:HB3	1:A:228:TRP:CZ3	2.47	0.50
1:B:167:ARG:O	1:B:170:LEU:HB3	2.12	0.50
1:B:10:ALA:HB2	1:B:25:CYS:HB3	1.94	0.50
1:B:26:MET:HB3	1:B:44:LEU:HD13	1.93	0.50
1:A:64:SER:O	1:A:65:ILE:C	2.55	0.50
1:A:109:ALA:HB1	1:A:115:LYS:HG3	1.94	0.50
1:A:192:ALA:O	1:A:196:PHE:HB3	2.12	0.50
1:A:10:ALA:HB2	1:A:25:CYS:HB3	1.94	0.50
1:A:126:TYR:CD2	1:A:144:GLN:HB3	2.46	0.50
1:B:98:LEU:HD11	1:B:126:TYR:CD1	2.46	0.50
1:A:32:GLN:O	1:A:104:PHE:HE1	1.95	0.49
1:A:3:LYS:O	1:A:7:VAL:HG23	2.12	0.49
1:A:17:GLU:O	1:A:17:GLU:HG2	2.12	0.49
1:B:14:GLU:C	1:B:16:ALA:N	2.71	0.49
1:B:98:LEU:HD21	1:B:126:TYR:CD1	2.48	0.49
1:B:126:TYR:CD2	1:B:144:GLN:HB3	2.46	0.49
1:A:174:PHE:CD1	1:A:178:TYR:CE1	3.00	0.49
1:B:32:GLN:O	1:B:104:PHE:HE1	1.95	0.49
1:A:156:SER:HB2	1:A:167:ARG:CG	2.43	0.49
1:B:3:LYS:O	1:B:7:VAL:HG23	2.12	0.49
1:A:90:LEU:CD2	1:A:132:VAL:HG21	2.42	0.49
1:A:126:TYR:O	1:A:130:ALA:N	2.46	0.49
1:B:192:ALA:O	1:B:196:PHE:HB3	2.12	0.49
1:B:64:SER:O	1:B:65:ILE:C	2.55	0.49
1:B:90:LEU:CD2	1:B:132:VAL:HG21	2.42	0.49
1:A:193:LYS:HA	1:A:196:PHE:CD2	2.48	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:GLY:O	1:B:56:ARG:N	2.46	0.48
1:A:48:TYR:CD2	1:A:97:VAL:HB	2.48	0.48
1:A:53:GLY:O	1:A:56:ARG:N	2.46	0.48
1:A:193:LYS:O	1:A:197:ASP:HB2	2.14	0.48
1:A:197:ASP:HA	1:A:200:ILE:CD1	2.44	0.48
1:B:156:SER:HB2	1:B:167:ARG:CG	2.43	0.48
1:B:48:TYR:CD2	1:B:97:VAL:HB	2.48	0.48
1:B:17:GLU:O	1:B:17:GLU:HG2	2.12	0.48
1:B:173:ASN:HA	1:B:176:VAL:HG12	1.95	0.48
1:A:14:GLU:C	1:A:16:ALA:N	2.71	0.48
1:A:173:ASN:HA	1:A:176:VAL:HG12	1.95	0.48
1:B:193:LYS:O	1:B:197:ASP:HB2	2.14	0.48
1:A:98:LEU:HD21	1:A:126:TYR:CD1	2.48	0.47
1:B:193:LYS:HA	1:B:196:PHE:CD2	2.48	0.47
1:B:151:GLU:O	1:B:155:ILE:HG13	2.14	0.47
1:A:9:LYS:HD2	1:A:25:CYS:SG	2.55	0.47
1:A:53:GLY:O	1:A:54:ALA:C	2.57	0.47
1:A:145:SER:O	1:A:148:ALA:HB3	2.15	0.47
1:B:174:PHE:CD1	1:B:178:TYR:CE1	3.00	0.47
1:A:26:MET:CB	1:A:44:LEU:HD12	2.44	0.47
1:A:103:LYS:O	1:A:104:PHE:CD2	2.67	0.47
1:A:151:GLU:O	1:A:155:ILE:HG13	2.14	0.47
1:B:26:MET:CB	1:B:44:LEU:HD12	2.44	0.47
1:B:221:LEU:HD23	1:B:221:LEU:N	2.28	0.47
1:B:61:VAL:HG22	1:B:62:VAL:N	2.29	0.47
1:B:103:LYS:O	1:B:104:PHE:CD2	2.67	0.47
1:B:145:SER:O	1:B:148:ALA:HB3	2.15	0.47
1:A:117:PHE:CD1	1:A:117:PHE:C	2.90	0.47
1:A:189:CYS:O	1:A:192:ALA:HB3	2.15	0.47
1:B:19:TYR:CD2	1:B:50:ASN:ND2	2.83	0.47
1:A:19:TYR:CD2	1:A:50:ASN:ND2	2.83	0.47
1:A:61:VAL:HG22	1:A:62:VAL:N	2.29	0.47
1:A:137:ASP:O	1:A:141:ILE:HG13	2.15	0.47
1:A:170:LEU:C	1:A:170:LEU:HD12	2.40	0.47
1:B:99:SER:O	1:B:100:LEU:C	2.58	0.46
1:A:37:SER:HB3	1:A:40:GLU:CD	2.40	0.46
1:B:9:LYS:HD2	1:B:25:CYS:SG	2.55	0.46
1:B:37:SER:HB3	1:B:40:GLU:CD	2.40	0.46
1:B:197:ASP:HA	1:B:200:ILE:CD1	2.44	0.46
1:A:46:VAL:HG23	1:A:49:LYS:HE3	1.97	0.46
1:A:166:ILE:HG12	1:A:167:ARG:N	2.31	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:SER:O	1:A:100:LEU:C	2.58	0.46
1:B:46:VAL:HG23	1:B:49:LYS:HE3	1.97	0.46
1:B:53:GLY:O	1:B:54:ALA:C	2.57	0.46
1:B:137:ASP:O	1:B:141:ILE:HG13	2.15	0.46
1:B:170:LEU:HD12	1:B:170:LEU:C	2.40	0.46
1:B:117:PHE:CD1	1:B:117:PHE:C	2.90	0.46
1:B:189:CYS:O	1:B:192:ALA:HB3	2.15	0.46
1:A:64:SER:O	1:A:66:GLU:N	2.50	0.45
1:B:43:LEU:H	1:B:43:LEU:CD1	2.29	0.45
1:A:46:VAL:O	1:A:47:ALA:C	2.59	0.45
1:A:196:PHE:O	1:A:199:ALA:HB3	2.16	0.45
1:B:46:VAL:O	1:B:47:ALA:C	2.59	0.45
1:B:143:ASP:O	1:B:147:GLN:HG3	2.16	0.45
1:A:52:VAL:HG11	1:A:125:TYR:CE1	2.50	0.45
1:B:43:LEU:CD1	1:B:43:LEU:N	2.80	0.45
1:B:90:LEU:HD21	1:B:132:VAL:CG2	2.47	0.45
1:B:196:PHE:O	1:B:199:ALA:HB3	2.16	0.45
1:A:43:LEU:H	1:A:43:LEU:CD1	2.29	0.45
1:B:126:TYR:O	1:B:130:ALA:N	2.46	0.45
1:A:54:ALA:O	1:A:57:SER:HB3	2.17	0.44
1:A:218:MET:O	1:A:219:GLN:C	2.60	0.44
1:B:142:VAL:HG13	1:B:143:ASP:N	2.32	0.44
1:B:166:ILE:HG12	1:B:167:ARG:N	2.31	0.44
1:A:6:LEU:HD13	1:A:28:SER:OG	2.18	0.44
1:A:119:LEU:HD23	1:A:151:GLU:HB3	2.00	0.44
1:A:142:VAL:HG13	1:A:143:ASP:N	2.32	0.44
1:A:101:LEU:HA	1:A:101:LEU:HD12	1.87	0.44
1:A:42:ASN:O	1:A:45:SER:OG	2.36	0.44
1:A:90:LEU:HD21	1:A:132:VAL:CG2	2.47	0.44
1:A:43:LEU:CD1	1:A:43:LEU:N	2.80	0.44
1:B:64:SER:O	1:B:66:GLU:N	2.50	0.44
1:B:119:LEU:HD22	1:B:155:ILE:HD12	1.99	0.44
1:B:42:ASN:O	1:B:45:SER:OG	2.36	0.44
1:B:218:MET:O	1:B:219:GLN:C	2.60	0.44
1:A:143:ASP:O	1:A:147:GLN:HG3	2.16	0.43
1:A:196:PHE:CD2	1:A:197:ASP:N	2.86	0.43
1:B:6:LEU:HD13	1:B:28:SER:OG	2.18	0.43
1:B:87:GLU:HG3	1:B:132:VAL:HG11	1.99	0.43
1:B:196:PHE:CD2	1:B:197:ASP:N	2.86	0.43
1:A:38:ASN:HA	1:A:41:ARG:HB3	2.00	0.43
1:A:89:GLU:H	1:A:89:GLU:HG2	1.66	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:LEU:HD22	1:A:155:ILE:HD12	1.99	0.43
1:B:36:LEU:O	1:B:37:SER:C	2.61	0.43
1:B:52:VAL:HG11	1:B:125:TYR:CE1	2.50	0.43
1:A:221:LEU:HD23	1:A:221:LEU:N	2.28	0.43
1:B:142:VAL:CG1	1:B:143:ASP:N	2.82	0.43
1:B:54:ALA:O	1:B:57:SER:HB3	2.17	0.43
1:B:41:ARG:HE	1:B:117:PHE:HD2	1.67	0.43
1:B:119:LEU:HD23	1:B:151:GLU:HB3	1.99	0.43
1:A:8:GLN:N	1:A:8:GLN:OE1	2.52	0.43
1:B:37:SER:HB3	1:B:40:GLU:HG3	2.01	0.43
1:B:86:ILE:HA	1:B:89:GLU:HG2	2.01	0.42
1:B:96:ASP:O	1:B:100:LEU:HD13	2.19	0.42
1:A:36:LEU:O	1:A:37:SER:C	2.61	0.42
1:A:37:SER:HB3	1:A:40:GLU:HG3	2.01	0.42
1:A:41:ARG:HE	1:A:117:PHE:HD2	1.67	0.42
1:A:142:VAL:CG1	1:A:143:ASP:N	2.82	0.42
1:B:38:ASN:HA	1:B:41:ARG:HB3	2.00	0.42
1:B:109:ALA:HB3	1:B:115:LYS:CG	2.49	0.42
1:B:8:GLN:N	1:B:8:GLN:OE1	2.52	0.42
1:B:218:MET:HE2	1:B:218:MET:HB3	1.90	0.42
1:A:86:ILE:C	1:A:88:THR:N	2.75	0.42
1:A:107:PRO:O	1:A:108:ASN:HB2	2.19	0.42
1:A:43:LEU:HD13	1:A:43:LEU:H	1.85	0.42
1:A:83:ARG:O	1:A:86:ILE:HG13	2.20	0.42
1:A:179:TYR:HD2	1:A:228:TRP:HE1	1.62	0.42
1:A:86:ILE:HA	1:A:89:GLU:HG2	2.01	0.42
1:A:87:GLU:HG3	1:A:132:VAL:HG11	1.99	0.42
1:A:87:GLU:HG2	1:A:91:ARG:HH21	1.83	0.42
1:A:156:SER:CB	1:A:167:ARG:HG3	2.49	0.42
1:B:43:LEU:HD13	1:B:43:LEU:H	1.85	0.42
1:B:83:ARG:O	1:B:86:ILE:HG13	2.20	0.42
1:B:107:PRO:O	1:B:108:ASN:HB2	2.19	0.42
1:A:129:LEU:HD12	1:A:129:LEU:HA	1.78	0.42
1:B:89:GLU:HG2	1:B:89:GLU:H	1.66	0.42
1:B:46:VAL:O	1:B:49:LYS:N	2.53	0.41
1:B:82:TYR:O	1:B:86:ILE:HG12	2.20	0.41
1:A:82:TYR:O	1:A:86:ILE:HG12	2.20	0.41
1:A:52:VAL:HG23	1:A:90:LEU:HD12	2.02	0.41
1:B:26:MET:O	1:B:29:VAL:N	2.54	0.41
1:A:26:MET:O	1:A:29:VAL:N	2.54	0.41
1:A:96:ASP:O	1:A:100:LEU:HD13	2.19	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:ALA:HB3	1:A:115:LYS:CG	2.50	0.41
1:A:149:TYR:O	1:A:152:ALA:N	2.54	0.41
1:A:32:GLN:CA	1:A:32:GLN:HE21	2.33	0.41
1:B:170:LEU:O	1:B:171:ALA:C	2.64	0.41
1:A:46:VAL:O	1:A:49:LYS:N	2.53	0.41
1:B:52:VAL:HG23	1:B:90:LEU:HD12	2.02	0.41
1:A:26:MET:HB3	1:A:44:LEU:HD12	2.02	0.41
1:B:30:THR:HG21	1:B:100:LEU:CD2	2.47	0.41
1:B:32:GLN:CA	1:B:32:GLN:HE21	2.33	0.41
1:B:51:VAL:HG12	1:B:93:ILE:HG21	2.03	0.41
1:A:124:ASP:OD1	1:A:149:TYR:OH	2.38	0.40
1:A:179:TYR:CD2	1:A:228:TRP:NE1	2.83	0.40
1:B:26:MET:HB3	1:B:44:LEU:HD12	2.02	0.40
1:B:87:GLU:HG2	1:B:91:ARG:HH21	1.83	0.40
1:B:156:SER:CB	1:B:167:ARG:HG3	2.49	0.40
1:B:58:SER:O	1:B:62:VAL:HG12	2.21	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	183/245 (75%)	131 (72%)	41 (22%)	11 (6%)	1	13
1	B	183/245 (75%)	131 (72%)	41 (22%)	11 (6%)	1	13
2	P	4/15 (27%)	0	0	4 (100%)	0	0
2	Q	4/15 (27%)	0	0	4 (100%)	0	0
All	All	374/520 (72%)	262 (70%)	82 (22%)	30 (8%)	1	8

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	ALA
1	A	108	ASN
2	P	257	SER
2	P	260	THR
2	P	261	PRO
1	B	34	ALA
1	B	108	ASN
2	Q	257	SER
2	Q	260	THR
2	Q	261	PRO
1	A	5	GLU
1	A	51	VAL
1	A	66	GLU
1	A	104	PHE
1	B	5	GLU
1	B	51	VAL
1	B	66	GLU
1	B	104	PHE
1	A	133	ALA
2	P	258	THR
1	B	133	ALA
2	Q	258	THR
1	A	100	LEU
1	B	100	LEU
1	A	2	ASP
1	A	65	ILE
1	A	99	SER
1	B	2	ASP
1	B	65	ILE
1	B	99	SER

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	166/209 (79%)	123 (74%)	43 (26%)	0 4
1	B	166/209 (79%)	123 (74%)	43 (26%)	0 4

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	332/418 (79%)	246 (74%)	86 (26%)	0 4

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	8	GLN
1	A	9	LYS
1	A	17	GLU
1	A	18	ARG
1	A	22	MET
1	A	30	THR
1	A	32	GLN
1	A	36	LEU
1	A	42	ASN
1	A	45	SER
1	A	56	ARG
1	A	61	VAL
1	A	62	VAL
1	A	67	GLN
1	A	76	GLN
1	A	78	MET
1	A	81	GLU
1	A	86	ILE
1	A	89	GLU
1	A	91	ARG
1	A	92	ASP
1	A	93	ILE
1	A	99	SER
1	A	101	LEU
1	A	105	LEU
1	A	108	ASN
1	A	116	VAL
1	A	129	LEU
1	A	132	VAL
1	A	144	GLN
1	A	149	TYR
1	A	166	ILE
1	A	168	LEU
1	A	170	LEU
1	A	189	CYS
1	A	196	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	213	ASP
1	A	214	SER
1	A	219	GLN
1	A	221	LEU
1	A	222	ARG
1	A	224	ASN
1	B	1	MET
1	B	8	GLN
1	B	9	LYS
1	B	17	GLU
1	B	18	ARG
1	B	22	MET
1	B	30	THR
1	B	32	GLN
1	B	36	LEU
1	B	42	ASN
1	B	45	SER
1	B	56	ARG
1	B	61	VAL
1	B	62	VAL
1	B	67	GLN
1	B	76	GLN
1	B	78	MET
1	B	81	GLU
1	B	86	ILE
1	B	89	GLU
1	B	91	ARG
1	B	92	ASP
1	B	93	ILE
1	B	99	SER
1	B	101	LEU
1	B	105	LEU
1	B	108	ASN
1	B	116	VAL
1	B	129	LEU
1	B	132	VAL
1	B	144	GLN
1	B	149	TYR
1	B	166	ILE
1	B	168	LEU
1	B	170	LEU
1	B	189	CYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	196	PHE
1	B	213	ASP
1	B	214	SER
1	B	219	GLN
1	B	221	LEU
1	B	222	ARG
1	B	224	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	42	ASN
1	A	67	GLN
1	A	144	GLN
1	B	32	GLN
1	B	42	ASN
1	B	67	GLN
1	B	144	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	P	259	2	8,9,10	2.61	2 (25%)	7,12,14	9.02	3 (42%)
2	SEP	Q	259	2	8,9,10	2.60	2 (25%)	7,12,14	9.03	3 (42%)

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	P	259	2	-	5/6/8/10	-
2	SEP	Q	259	2	-	5/6/8/10	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	259	SEP	OG-CB	-5.20	1.24	1.44
2	Q	259	SEP	OG-CB	-5.18	1.25	1.44
2	P	259	SEP	P-OG	4.60	1.74	1.60
2	Q	259	SEP	P-OG	4.59	1.74	1.60

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	259	SEP	OG-CB-CA	23.64	131.15	108.14
2	P	259	SEP	OG-CB-CA	23.62	131.13	108.14
2	Q	259	SEP	O2P-P-OG	-2.23	100.87	106.67
2	P	259	SEP	O2P-P-OG	-2.22	100.88	106.67
2	Q	259	SEP	O3P-P-O2P	2.06	115.53	107.80
2	P	259	SEP	O3P-P-O2P	2.05	115.49	107.80

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	P	259	SEP	N-CA-CB-OG
2	P	259	SEP	C-CA-CB-OG
2	P	259	SEP	CB-OG-P-O1P
2	P	259	SEP	CB-OG-P-O2P
2	P	259	SEP	CB-OG-P-O3P
2	Q	259	SEP	N-CA-CB-OG
2	Q	259	SEP	C-CA-CB-OG
2	Q	259	SEP	CB-OG-P-O1P
2	Q	259	SEP	CB-OG-P-O2P
2	Q	259	SEP	CB-OG-P-O3P

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	259	SEP	1	0
2	Q	259	SEP	1	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	A	195/245 (79%)	2.01	90 (46%)  	32, 77, 99, 100	0
1	B	195/245 (79%)	2.04	95 (48%)  	32, 77, 99, 100	0
2	P	5/15 (33%)	0.88	0  	45, 57, 62, 65	0
2	Q	5/15 (33%)	2.00	3 (60%)  	45, 57, 62, 65	0
All	All	400/520 (76%)	2.01	188 (47%)  	32, 77, 99, 100	0

All (188) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	145	SER	6.3
1	B	99	SER	5.3
1	B	38	ASN	5.0
1	B	33	GLY	4.8
1	B	6	LEU	4.7
1	B	3	LYS	4.6
1	A	17	GLU	4.5
1	B	41	ARG	4.4
1	A	168	LEU	4.2
1	B	169	GLY	4.1
1	A	123	GLY	4.1
1	A	84	GLU	4.1
1	B	49	LYS	4.1
1	B	75	LYS	4.0
1	B	22	MET	4.0
1	A	157	LYS	3.9
1	A	3	LYS	3.9
1	B	34	ALA	3.9
1	B	223	ASP	3.8
1	A	198	GLU	3.8
1	B	168	LEU	3.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	7	VAL	3.7
1	A	169	GLY	3.7
1	A	217	ILE	3.7
1	A	56	ARG	3.6
1	B	17	GLU	3.6
1	B	67	GLN	3.6
1	B	137	ASP	3.6
1	A	201	ALA	3.6
1	A	33	GLY	3.6
1	A	6	LEU	3.5
1	A	81	GLU	3.5
1	B	149	TYR	3.4
1	A	148	ALA	3.4
1	B	1	MET	3.4
1	B	191	LEU	3.4
1	A	22	MET	3.4
1	A	196	PHE	3.3
1	B	60	ARG	3.3
1	B	153	PHE	3.3
1	B	226	THR	3.3
1	B	150	GLN	3.3
1	B	107	PRO	3.3
1	A	166	ILE	3.3
1	B	123	GLY	3.3
1	A	175	SER	3.3
1	B	76	GLN	3.3
1	A	52	VAL	3.2
1	A	177	PHE	3.2
1	B	2	ASP	3.2
1	A	44	LEU	3.2
1	A	2	ASP	3.2
1	A	10	ALA	3.2
1	B	15	GLN	3.1
1	A	223	ASP	3.1
1	B	152	ALA	3.1
1	A	29	VAL	3.0
1	B	56	ARG	3.0
1	B	138	LYS	3.0
1	B	179	TYR	3.0
1	B	134	ALA	3.0
1	B	146	GLN	3.0
1	A	26	MET	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	219	GLN	2.9
1	A	75	LYS	2.9
1	A	218	MET	2.9
1	B	145	SER	2.9
1	A	105	LEU	2.9
1	A	135	GLY	2.9
1	B	147	GLN	2.9
1	A	60	ARG	2.9
2	Q	256	ARG	2.9
1	B	142	VAL	2.9
1	A	49	LYS	2.9
1	B	12	LEU	2.9
1	A	174	PHE	2.9
1	B	112	ALA	2.8
1	A	112	ALA	2.8
1	A	5	GLU	2.8
1	A	154	GLU	2.8
1	A	79	ALA	2.8
1	B	143	ASP	2.8
1	A	176	VAL	2.8
1	A	19	TYR	2.8
1	B	113	GLU	2.8
1	B	198	GLU	2.8
1	A	153	PHE	2.8
1	A	195	ALA	2.8
1	B	35	GLU	2.7
1	A	11	LYS	2.7
1	B	196	PHE	2.7
1	A	80	ARG	2.7
1	B	9	LYS	2.7
2	Q	261	PRO	2.7
1	B	81	GLU	2.7
1	A	67	GLN	2.7
1	A	34	ALA	2.7
1	A	113	GLU	2.7
1	A	228	TRP	2.7
1	B	77	GLN	2.7
1	A	73	GLU	2.7
1	B	61	VAL	2.6
1	A	77	GLN	2.6
1	A	85	LYS	2.6
1	B	171	ALA	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	139	LYS	2.6
1	A	96	ASP	2.6
1	A	106	ILE	2.6
1	A	155	ILE	2.6
1	B	4	ASN	2.6
1	A	167	ARG	2.6
1	B	227	LEU	2.6
1	A	179	TYR	2.5
1	A	200	ILE	2.5
1	B	175	SER	2.5
1	A	103	LYS	2.5
1	A	32	GLN	2.5
1	A	74	LYS	2.5
1	A	136	ASP	2.5
1	A	156	SER	2.5
1	B	27	LYS	2.5
1	B	141	ILE	2.5
1	B	7	VAL	2.5
1	B	128	TYR	2.5
1	A	24	ALA	2.4
1	A	222	ARG	2.4
1	A	214	SER	2.4
1	A	150	GLN	2.4
1	B	132	VAL	2.4
1	B	166	ILE	2.4
1	A	149	TYR	2.4
1	B	216	LEU	2.4
1	B	80	ARG	2.4
1	A	100	LEU	2.4
1	A	128	TYR	2.4
1	A	197	ASP	2.4
1	A	1	MET	2.4
1	B	199	ALA	2.3
1	A	15	GLN	2.3
1	A	63	SER	2.3
1	A	221	LEU	2.3
1	B	105	LEU	2.3
1	A	99	SER	2.3
1	B	47	ALA	2.3
1	B	88	THR	2.3
1	A	36	LEU	2.3
1	B	106	ILE	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	174	PHE	2.3
1	B	26	MET	2.3
2	Q	258	THR	2.2
1	B	173	ASN	2.2
1	B	225	LEU	2.2
1	A	65	ILE	2.2
1	B	135	GLY	2.2
1	B	45	SER	2.2
1	A	141	ILE	2.2
1	B	115	LYS	2.2
1	A	23	ALA	2.2
1	B	124	ASP	2.2
1	B	36	LEU	2.2
1	B	170	LEU	2.2
1	B	28	SER	2.1
1	A	116	VAL	2.1
1	B	97	VAL	2.1
1	A	133	ALA	2.1
1	B	201	ALA	2.1
1	B	32	GLN	2.1
1	A	107	PRO	2.1
1	B	83	ARG	2.1
1	B	24	ALA	2.1
1	A	38	ASN	2.1
1	A	132	VAL	2.1
1	A	40	GLU	2.1
1	A	9	LYS	2.1
1	B	157	LYS	2.1
1	B	217	ILE	2.1
1	A	27	LYS	2.0
1	B	133	ALA	2.0
1	B	31	GLU	2.0
1	B	63	SER	2.0
1	B	130	ALA	2.0
1	B	222	ARG	2.0
1	B	66	GLU	2.0
1	B	131	GLU	2.0
1	B	23	ALA	2.0
1	B	109	ALA	2.0
1	B	156	SER	2.0
1	B	188	ALA	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(Å ²)	Q<0.9
2	SEP	Q	259	10/11	0.73	0.17	46,66,66,66	0
2	SEP	P	259	10/11	0.75	0.20	46,66,66,66	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.