



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:03 PM UTC

PDB ID : 4ZHE / pdb_00004zhe
Title : Crystal structure of the SeMet substituted Topless related protein 2 (TPR2)
N-terminal domain (1-209) from rice
Authors : Ke, J.; Ma, H.; Gu, X.; Brunzelle, J.S.; Xu, H.E.; Melcher, K.
Deposited on : 2015-04-24
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

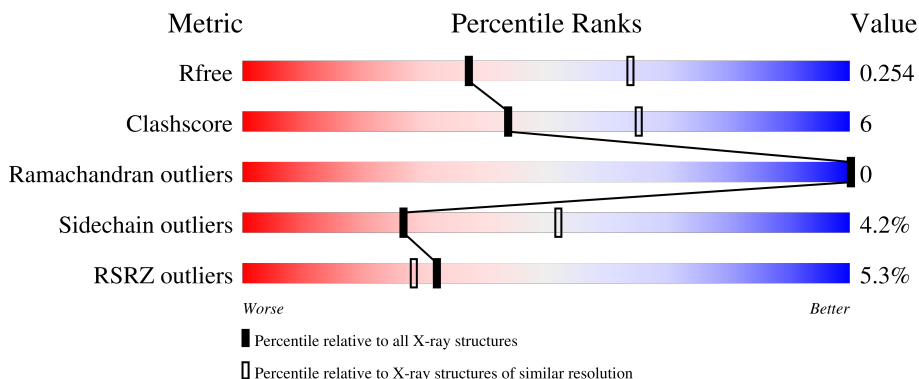
1 Overall quality at a glance

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))
R_{free}	180053	5829 (2.50-2.50)
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	A	209	4% 80% 16% ..
1	B	209	3% 80% 15% ..
1	C	209	5% 77% 16% • 5%
1	D	209	8% 79% 17% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7182 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 ASPR2 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	Se	0	0	0
			1716	1107	287	316	2	4			
1	B	203	Total	C	N	O	S	Se	0	0	0
			1708	1102	286	314	2	4			
1	C	199	Total	C	N	O	S	Se	0	0	0
			1675	1081	279	309	2	4			
1	D	206	Total	C	N	O	S	Se	0	0	0
			1705	1100	282	318	1	4			

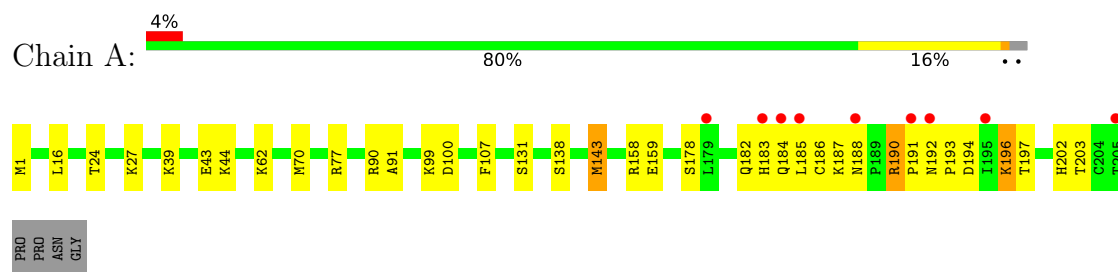
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	126	Total	O	0	0
			126	126		
2	B	134	Total	O	0	0
			134	134		
2	C	44	Total	O	0	0
			44	44		
2	D	74	Total	O	0	0
			74	74		

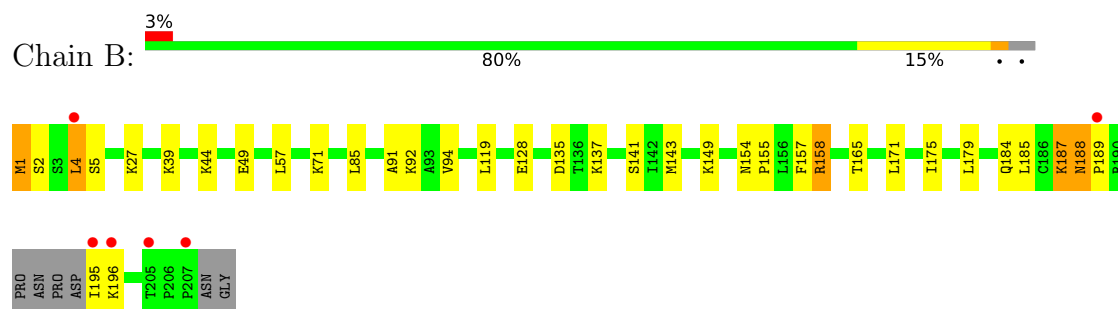
3 Residue-property plots [i](#)

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.

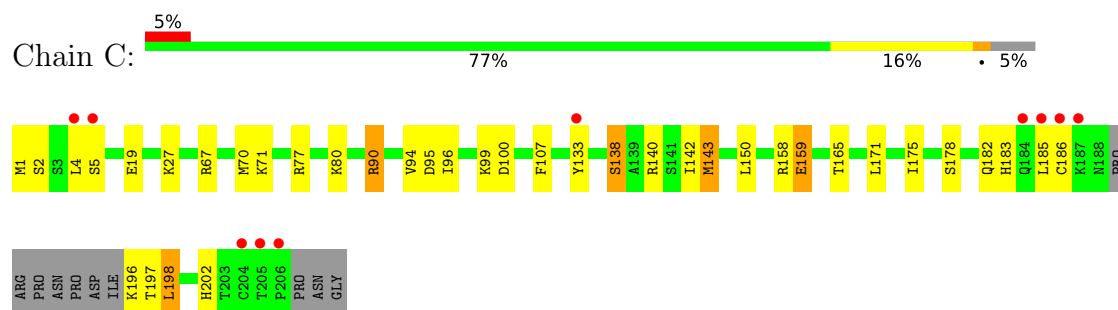
- Molecule 1: ASPR2 protein



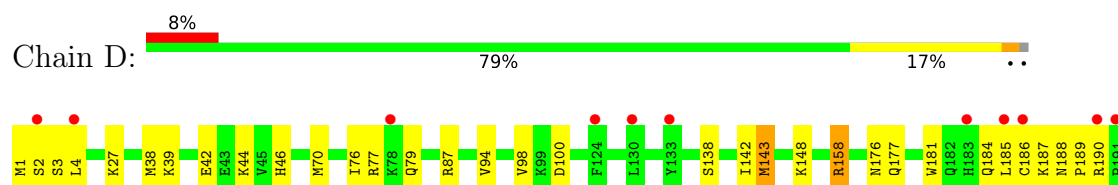
- Molecule 1: ASPR2 protein

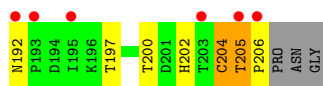


- Molecule 1: ASPR2 protein



- Molecule 1: ASPR2 protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.15Å 111.88Å 144.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.28 – 2.50 44.28 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.28-2.50) 100.0 (44.28-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.199 , 0.242 0.214 , 0.254	Depositor DCC
R_{free} test set	2038 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	48.8	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7182	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.52% 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

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	0.66	1/1748 (0.1%)	0.95	0/2343
1	B	0.71	4/1739 (0.2%)	1.00	8/2329 (0.3%)
1	C	0.68	4/1704 (0.2%)	1.01	8/2279 (0.4%)
1	D	0.83	6/1737 (0.3%)	1.10	9/2333 (0.4%)
All	All	0.73	15/6928 (0.2%)	1.02	25/9284 (0.3%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	186	CYS	CA-CB	8.74	1.66	1.53
1	D	185	LEU	CA-CB	8.53	1.68	1.53
1	B	128	GLU	CG-CD	-7.95	1.32	1.52
1	D	186	CYS	N-CA	7.87	1.54	1.46
1	C	67	ARG	CZ-NH2	-7.22	1.24	1.33
1	B	128	GLU	CB-CG	-7.21	1.30	1.52
1	C	67	ARG	CD-NE	-6.87	1.36	1.46
1	D	185	LEU	CA-C	5.60	1.60	1.52
1	B	128	GLU	CD-OE1	-5.55	1.14	1.25
1	C	159	GLU	CG-CD	-5.53	1.38	1.52
1	D	176	ASN	CG-OD1	-5.50	1.13	1.23
1	B	149	LYS	CE-NZ	-5.35	1.33	1.49
1	C	67	ARG	CZ-NH1	-5.32	1.25	1.32
1	D	184	GLN	CA-CB	5.20	1.61	1.53
1	A	196	LYS	CD-CE	-5.04	1.37	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	90	ARG	NE-CZ-NH2	13.64	131.48	119.20
1	C	90	ARG	NE-CZ-NH1	-11.88	109.62	121.50
1	D	190	ARG	CA-C-N	-10.68	108.82	120.14
1	D	190	ARG	C-N-CA	-10.68	108.82	120.14
1	D	2	SER	N-CA-C	-10.55	101.11	112.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	157	PHE	N-CA-C	8.51	123.97	113.50
1	B	196	LYS	N-CA-C	8.16	121.56	110.55
1	B	187	LYS	N-CA-C	-8.03	97.91	110.36
1	D	184	GLN	N-CA-C	7.84	120.72	111.71
1	C	90	ARG	CD-NE-CZ	-7.67	113.67	124.40
1	D	181	TRP	N-CA-C	-7.36	104.28	113.18
1	C	185	LEU	N-CA-C	-7.24	94.79	108.17
1	D	192	ASN	CA-C-N	-7.07	111.00	119.84
1	D	192	ASN	C-N-CA	-7.07	111.00	119.84
1	B	185	LEU	CA-C-N	-6.32	114.07	122.85
1	B	185	LEU	C-N-CA	-6.32	114.07	122.85
1	D	186	CYS	N-CA-CB	6.25	123.44	111.00
1	B	128	GLU	N-CA-C	5.89	118.46	111.33
1	B	128	GLU	CA-CB-CG	-5.63	102.85	114.10
1	D	158	ARG	N-CA-C	5.40	116.86	110.97
1	C	90	ARG	CG-CD-NE	-5.24	100.47	112.00
1	C	67	ARG	CG-CD-NE	-5.19	100.58	112.00
1	C	90	ARG	CB-CG-CD	-5.10	99.58	111.30
1	C	197	THR	N-CA-C	5.06	116.69	109.14
1	B	184	GLN	N-CA-C	5.04	116.77	111.28

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	1716	0	1718	24	1
1	B	1708	0	1720	24	1
1	C	1675	0	1689	23	0
1	D	1705	0	1691	25	0
2	A	126	0	0	0	0
2	B	134	0	0	5	0
2	C	44	0	0	1	0
2	D	74	0	0	3	0
All	All	7182	0	6818	83	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 6.

All (83) 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:C:90:ARG:HH12	1:D:98:VAL:HG21	1.27	0.98
1:A:27:LYS:HG3	1:D:27:LYS:HG3	1.75	0.69
1:A:186:CYS:SG	1:A:187:LYS:N	2.66	0.67
1:C:90:ARG:NH1	1:D:98:VAL:HG21	2.06	0.66
1:A:190:ARG:HH11	1:A:190:ARG:HB2	1.60	0.66
1:B:44:LYS:NZ	2:B:301:HOH:O	2.28	0.66
1:B:27:LYS:HG3	1:C:27:LYS:HG3	1.79	0.65
1:B:195:ILE:HD11	1:C:4:LEU:HD11	1.80	0.63
1:D:76:ILE:HA	1:D:143:MSE:CE	2.33	0.59
1:D:148:LYS:NZ	2:D:302:HOH:O	2.37	0.58
1:C:140:ARG:HA	1:C:143:MSE:HE2	1.88	0.56
1:A:190:ARG:HH11	1:A:190:ARG:CB	2.18	0.55
1:C:70:MSE:HE1	1:C:107:PHE:CZ	2.40	0.55
1:D:138:SER:O	1:D:142:ILE:HG12	2.07	0.55
1:C:95:ASP:OD1	1:C:99:LYS:HE3	2.08	0.53
1:C:196:LYS:HG2	1:C:202:HIS:CE1	2.43	0.53
1:D:205:THR:O	1:D:205:THR:OG1	2.26	0.52
1:A:191:PRO:HA	1:A:192:ASN:C	2.34	0.52
1:D:1:MSE:C	1:D:3:SER:H	2.14	0.52
1:C:133:TYR:HB3	2:C:317:HOH:O	2.10	0.51
1:D:187:LYS:C	1:D:189:PRO:HD3	2.36	0.51
1:A:90:ARG:HD3	1:B:94:VAL:HG11	1.93	0.50
1:C:94:VAL:HG21	1:D:94:VAL:HG21	1.92	0.50
1:A:77:ARG:HD3	1:A:100:ASP:HB3	1.93	0.50
1:A:196:LYS:HG2	1:A:197:THR:N	2.26	0.49
1:D:204:CYS:O	1:D:206:PRO:HD3	2.12	0.49
1:B:135:ASP:OD2	1:B:137:LYS:NZ	2.46	0.49
1:D:188:ASN:N	1:D:189:PRO:HD3	2.27	0.49
1:A:39:LYS:O	1:A:43:GLU:HG3	2.13	0.49
1:B:4:LEU:HD11	1:C:198:LEU:HD13	1.94	0.48
1:C:138:SER:O	1:C:142:ILE:HG12	2.13	0.48
1:C:80:LYS:HE3	1:C:96:ILE:HG12	1.95	0.48
1:B:155:PRO:HA	1:B:158:ARG:HD2	1.95	0.48
1:C:77:ARG:HD3	1:C:100:ASP:HB3	1.96	0.48
1:C:178:SER:O	1:C:182:GLN:HG3	2.14	0.47
1:C:71:LYS:HG2	1:C:150:LEU:HD21	1.96	0.47
1:A:91:ALA:HB2	1:B:91:ALA:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1:MSE:SE	1:D:3:SER:HB3	2.65	0.47
1:A:90:ARG:HD3	1:B:94:VAL:CG1	2.45	0.46
1:A:44:LYS:HA	1:A:44:LYS:HD3	1.66	0.46
1:A:143:MSE:HB3	1:A:143:MSE:HE2	1.65	0.46
1:D:87:ARG:HD2	2:D:307:HOH:O	2.16	0.46
1:D:1:MSE:HG3	1:D:3:SER:HB3	1.97	0.46
1:B:2:SER:HB3	1:B:5:SER:OG	2.17	0.45
1:B:154:ASN:O	1:B:158:ARG:HG3	2.16	0.45
1:A:99:LYS:HA	1:A:99:LYS:HD3	1.68	0.45
1:C:158:ARG:HG2	1:C:159:GLU:HG2	1.97	0.45
1:D:197:THR:HG23	1:D:200:THR:H	1.80	0.45
1:C:196:LYS:HG2	1:C:202:HIS:ND1	2.32	0.45
1:B:188:ASN:N	1:B:189:PRO:HD3	2.31	0.44
1:B:85:LEU:HD21	1:B:119:LEU:HD23	2.00	0.44
1:D:79:GLN:HB3	1:D:143:MSE:HE3	1.99	0.44
1:A:16:LEU:HD13	1:A:24:THR:HG22	1.99	0.44
1:B:49:GLU:HG3	2:B:301:HOH:O	2.17	0.44
1:D:46:HIS:HD2	2:D:339:HOH:O	2.01	0.44
1:A:178:SER:O	1:A:182:GLN:HG3	2.17	0.44
1:A:178:SER:HB2	1:D:177:GLN:HB2	2.00	0.43
1:D:76:ILE:HA	1:D:143:MSE:HE2	1.99	0.43
1:D:77:ARG:HD3	1:D:100:ASP:HB3	1.99	0.43
1:A:62:LYS:HD3	1:A:62:LYS:HA	1.76	0.43
1:B:171:LEU:O	1:B:175:ILE:HG12	2.19	0.43
1:A:70:MSE:HE1	1:A:107:PHE:CZ	2.54	0.43
1:A:158:ARG:O	1:A:159:GLU:HB2	2.19	0.42
1:A:183:HIS:O	1:A:186:CYS:HB3	2.19	0.42
1:B:4:LEU:HD13	1:B:4:LEU:HA	1.85	0.42
1:D:38:MSE:HE3	1:D:42:GLU:HG3	2.01	0.42
1:B:92:LYS:NZ	2:B:307:HOH:O	2.52	0.42
1:D:79:GLN:CB	1:D:143:MSE:HE3	2.49	0.42
1:C:183:HIS:O	1:C:186:CYS:HB2	2.20	0.42
1:A:192:ASN:N	1:A:193:PRO:HA	2.35	0.42
1:B:188:ASN:ND2	2:B:309:HOH:O	2.52	0.41
1:B:1:MSE:HB3	1:B:2:SER:H	1.34	0.41
1:B:1:MSE:HE2	1:C:19:GLU:OE1	2.20	0.41
1:B:143:MSE:HB3	1:B:143:MSE:HE2	1.81	0.41
1:D:44:LYS:HD3	1:D:44:LYS:HA	1.67	0.41
1:A:70:MSE:HE1	1:A:107:PHE:CE1	2.56	0.41
1:B:158:ARG:NH1	2:B:311:HOH:O	2.54	0.41
1:C:90:ARG:O	1:C:94:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:LYS:O	1:A:188:ASN:C	2.63	0.40
1:B:4:LEU:HD12	1:C:198:LEU:HD22	2.04	0.40
1:C:171:LEU:O	1:C:175:ILE:HG12	2.20	0.40
1:D:4:LEU:HD23	1:D:4:LEU:HA	1.74	0.40
1:B:179:LEU:HD23	1:B:179:LEU:HA	1.92	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:A:131:SER:O	1:B:141:SER:OG[4_566]	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	203/209 (97%)	198 (98%)	5 (2%)	0	100	100
1	B	199/209 (95%)	198 (100%)	1 (0%)	0	100	100
1	C	195/209 (93%)	194 (100%)	1 (0%)	0	100	100
1	D	204/209 (98%)	196 (96%)	8 (4%)	0	100	100
All	All	801/836 (96%)	786 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	190/193 (98%)	181 (95%)	9 (5%)	23	47
1	B	191/193 (99%)	182 (95%)	9 (5%)	23	47
1	C	187/193 (97%)	180 (96%)	7 (4%)	30	57
1	D	187/193 (97%)	180 (96%)	7 (4%)	30	57
All	All	755/772 (98%)	723 (96%)	32 (4%)	26	52

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	138	SER
1	A	143	MSE
1	A	184	GLN
1	A	185	LEU
1	A	190	ARG
1	A	194	ASP
1	A	202	HIS
1	A	203	THR
1	B	1	MSE
1	B	4	LEU
1	B	39	LYS
1	B	57	LEU
1	B	71	LYS
1	B	158	ARG
1	B	165	THR
1	B	187	LYS
1	B	188	ASN
1	C	1	MSE
1	C	2	SER
1	C	5	SER
1	C	138	SER
1	C	143	MSE
1	C	165	THR
1	C	198	LEU
1	D	39	LYS
1	D	70	MSE
1	D	143	MSE
1	D	158	ARG
1	D	202	HIS

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Mol	Chain	Res	Type
1	D	204	CYS
1	D	205	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	HIS
1	A	180	ASN
1	B	127	ASN
1	C	176	ASN
1	C	183	HIS
1	C	202	HIS
1	D	30	GLN
1	D	117	GLN
1	D	176	ASN
1	D	202	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

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 ⓘ

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	201/209 (96%)	-0.04	9 (4%)	38 33	26, 42, 143, 166	0
1	B	199/209 (95%)	-0.05	6 (3%)	52 48	25, 45, 90, 127	0
1	C	195/209 (93%)	0.48	10 (5%)	33 29	43, 67, 107, 142	0
1	D	202/209 (96%)	0.36	17 (8%)	17 15	34, 57, 135, 182	0
All	All	797/836 (95%)	0.19	42 (5%)	32 28	25, 55, 119, 182	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	205	THR	5.3
1	B	195	ILE	4.7
1	C	185	LEU	4.4
1	C	204	CYS	3.7
1	B	196	LYS	3.7
1	C	206	PRO	3.6
1	D	205	THR	3.5
1	B	207	PRO	3.5
1	D	183	HIS	3.3
1	A	192	ASN	3.3
1	D	185	LEU	3.3
1	A	188	ASN	3.2
1	D	190	ARG	3.1
1	A	183	HIS	3.1
1	D	124	PHE	3.0
1	A	195	ILE	3.0
1	A	185	LEU	3.0
1	D	130	LEU	2.9
1	D	191	PRO	2.8
1	D	2	SER	2.8
1	C	205	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	186	CYS	2.8
1	D	206	PRO	2.8
1	C	187	LYS	2.8
1	D	195	ILE	2.7
1	A	191	PRO	2.7
1	D	78	LYS	2.5
1	D	203	THR	2.5
1	C	133	TYR	2.4
1	C	4	LEU	2.4
1	C	184	GLN	2.4
1	B	4	LEU	2.3
1	A	184	GLN	2.3
1	D	193	PRO	2.3
1	C	186	CYS	2.2
1	C	5	SER	2.2
1	B	189	PRO	2.2
1	A	179	LEU	2.1
1	B	205	THR	2.1
1	D	4	LEU	2.1
1	D	192	ASN	2.0
1	D	133	TYR	2.0

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

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

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.