



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:28 PM UTC

PDB ID : 4S2T / pdb_00004s2t
Title : Crystal structure of X-prolyl aminopeptidase from *Caenorhabditis elegans*: a cytosolic enzyme with a di-nuclear active site
Authors : Iyer, S.; La-Borde, P.; Payne, K.A.P.; Parsons, M.R.; Turner, A.J.; Isaac, R.E.; Acharya, K.R.
Deposited on : 2015-01-22
Resolution : 2.15 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

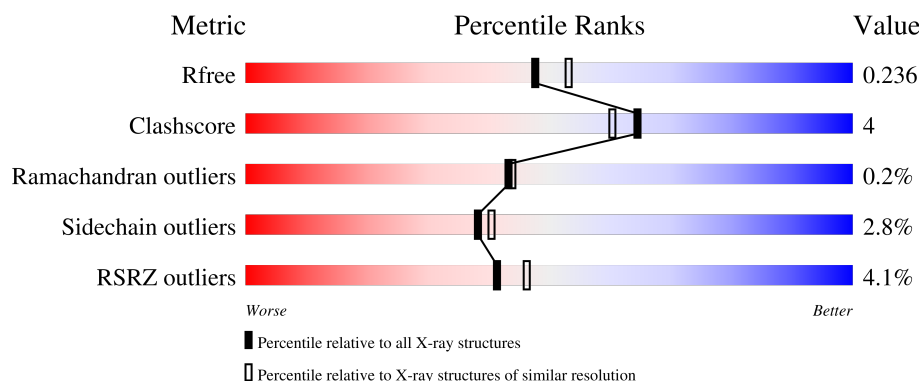
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	P	639	4% 88% 8% .
1	Q	639	3% 87% 8% ..
2	A	5	60% 40% 40% 20%
2	B	5	60% 40% 40% 20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	01B	A	1	-	-	X	-
2	01B	B	1	-	-	X	X
4	SO4	P	703	-	-	X	-
4	SO4	P	705	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10190 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 Protein APP-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	P	615	Total	C	N	O	S	25	0	0
			4876	3110	834	918	14			
1	Q	612	Total	C	N	O	S	13	0	0
			4859	3100	829	916	14			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	-22	MET	-	expression tag	UNP O44750
P	-21	GLY	-	expression tag	UNP O44750
P	-20	HIS	-	expression tag	UNP O44750
P	-19	HIS	-	expression tag	UNP O44750
P	-18	HIS	-	expression tag	UNP O44750
P	-17	HIS	-	expression tag	UNP O44750
P	-16	HIS	-	expression tag	UNP O44750
P	-15	HIS	-	expression tag	UNP O44750
P	-14	HIS	-	expression tag	UNP O44750
P	-13	HIS	-	expression tag	UNP O44750
P	-12	HIS	-	expression tag	UNP O44750
P	-11	HIS	-	expression tag	UNP O44750
P	-10	SER	-	expression tag	UNP O44750
P	-9	SER	-	expression tag	UNP O44750
P	-8	GLY	-	expression tag	UNP O44750
P	-7	HIS	-	expression tag	UNP O44750
P	-6	ILE	-	expression tag	UNP O44750
P	-5	ASP	-	expression tag	UNP O44750
P	-4	ASP	-	expression tag	UNP O44750
P	-3	ASP	-	expression tag	UNP O44750
P	-2	ASP	-	expression tag	UNP O44750
P	-1	LYS	-	expression tag	UNP O44750
P	0	HIS	-	expression tag	UNP O44750
Q	-22	MET	-	expression tag	UNP O44750
Q	-21	GLY	-	expression tag	UNP O44750

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Chain	Residue	Modelled	Actual	Comment	Reference
Q	-20	HIS	-	expression tag	UNP O44750
Q	-19	HIS	-	expression tag	UNP O44750
Q	-18	HIS	-	expression tag	UNP O44750
Q	-17	HIS	-	expression tag	UNP O44750
Q	-16	HIS	-	expression tag	UNP O44750
Q	-15	HIS	-	expression tag	UNP O44750
Q	-14	HIS	-	expression tag	UNP O44750
Q	-13	HIS	-	expression tag	UNP O44750
Q	-12	HIS	-	expression tag	UNP O44750
Q	-11	HIS	-	expression tag	UNP O44750
Q	-10	SER	-	expression tag	UNP O44750
Q	-9	SER	-	expression tag	UNP O44750
Q	-8	GLY	-	expression tag	UNP O44750
Q	-7	HIS	-	expression tag	UNP O44750
Q	-6	ILE	-	expression tag	UNP O44750
Q	-5	ASP	-	expression tag	UNP O44750
Q	-4	ASP	-	expression tag	UNP O44750
Q	-3	ASP	-	expression tag	UNP O44750
Q	-2	ASP	-	expression tag	UNP O44750
Q	-1	LYS	-	expression tag	UNP O44750
Q	0	HIS	-	expression tag	UNP O44750

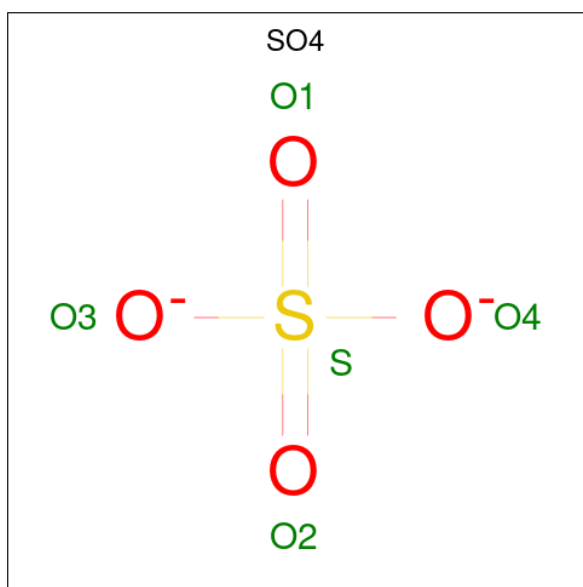
- Molecule 2 is a protein called apstatin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	A	4	Total	C	N	O	0	0	0
			32	23	4	5			
2	B	4	Total	C	N	O	0	0	0
			32	23	4	5			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	P	2	Total	Zn	0	0
			2	2		
3	Q	2	Total	Zn	0	0
			2	2		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	P	1	Total	O	S	0	0
			5	4	1		
4	P	1	Total	O	S	0	0
			5	4	1		
4	P	1	Total	O	S	0	0
			5	4	1		
4	Q	1	Total	O	S	0	0
			5	4	1		
4	Q	1	Total	O	S	0	0
			5	4	1		
4	Q	1	Total	O	S	0	0
			5	4	1		
4	Q	1	Total	O	S	0	0
			5	4	1		

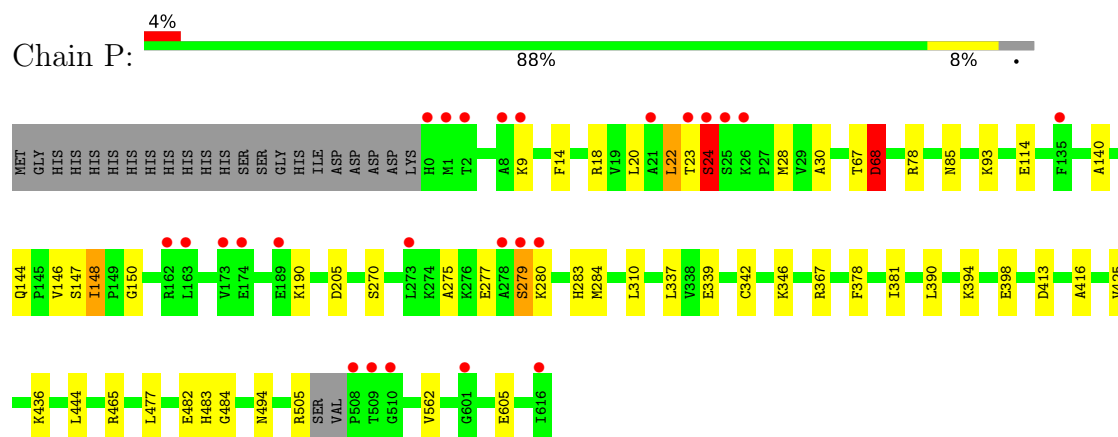
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	P	167	Total	O	0	0
			167	167		
5	Q	183	Total	O	0	0
			183	183		
5	A	1	Total	O	0	0
			1	1		
5	B	1	Total	O	0	0
			1	1		

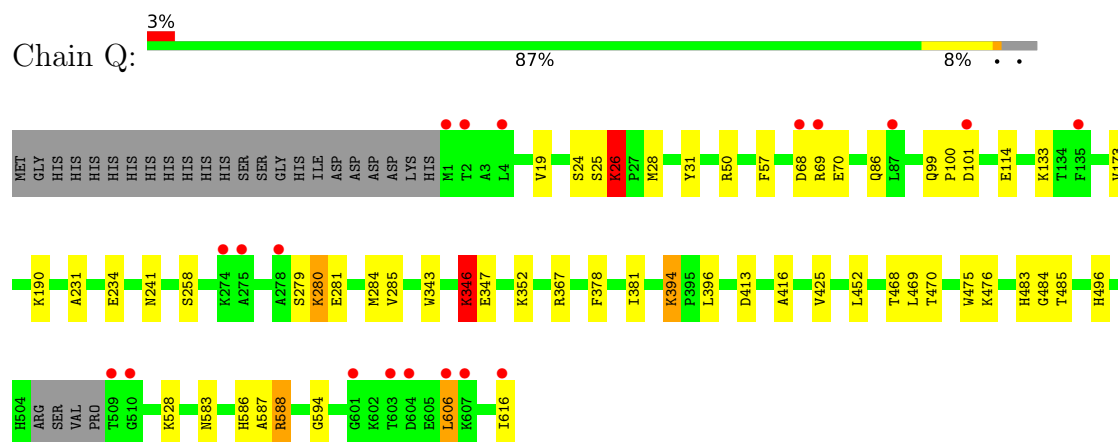
3 Residue-property plots [i](#)

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.

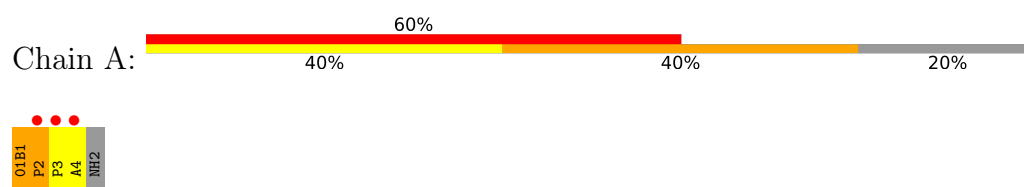
• Molecule 1: Protein APP-1



• Molecule 1: Protein APP-1



• Molecule 2: apstatin



• Molecule 2: apstatin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	140.59Å 86.81Å 113.13Å 90.00° 115.96° 90.00°	Depositor
Resolution (Å)	38.52 – 2.15 38.52 – 2.15	Depositor EDS
% Data completeness (in resolution range)	91.7 (38.52-2.15) 91.7 (38.52-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.205 , 0.247 (Not available) , 0.236	Depositor DCC
R_{free} test set	3099 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å ²)	32.8	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 27.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10190	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% 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: 01B, ZN, SO4

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	P	0.77	6/4990 (0.1%)	0.79	4/6765 (0.1%)
1	Q	0.76	10/4972 (0.2%)	0.80	7/6740 (0.1%)
2	A	2.57	2/20 (10.0%)	1.25	0/27
2	B	2.58	2/20 (10.0%)	1.25	0/27
All	All	0.78	20/10002 (0.2%)	0.80	11/13559 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	P	0	1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3	PRO	CA-C	-9.15	1.39	1.52
2	A	3	PRO	CA-C	-9.07	1.39	1.52
2	A	2	PRO	CA-C	-6.33	1.39	1.52
2	B	2	PRO	CA-C	-6.30	1.39	1.52
1	Q	26	LYS	C-N	6.12	1.40	1.33
1	Q	587	ALA	C-O	-5.74	1.17	1.24
1	Q	133	LYS	C-O	-5.60	1.17	1.24
1	Q	50	ARG	C-O	-5.41	1.17	1.24
1	Q	100	PRO	N-CD	5.35	1.55	1.47
1	Q	475	TRP	C-O	-5.34	1.17	1.24
1	P	148	ILE	C-N	5.32	1.40	1.33
1	P	342	CYS	C-O	-5.27	1.17	1.24
1	Q	588	ARG	C-O	-5.21	1.18	1.24
1	P	68	ASP	CA-C	-5.18	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	Q	586	HIS	C-O	-5.18	1.17	1.24
1	Q	476	LYS	C-O	-5.15	1.17	1.24
1	P	93	LYS	CG-CD	-5.12	1.37	1.52
1	Q	99	GLN	C-N	5.08	1.41	1.33
1	P	30	ALA	C-O	-5.04	1.17	1.23
1	P	436	LYS	C-O	-5.01	1.18	1.24

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	26	LYS	CA-C-N	-6.51	113.93	120.31
1	Q	26	LYS	C-N-CA	-6.51	113.93	120.31
1	Q	346	LYS	CB-CA-C	6.07	121.17	110.85
1	P	148	ILE	CA-C-N	-5.60	114.01	119.78
1	P	148	ILE	C-N-CA	-5.60	114.01	119.78
1	Q	99	GLN	CA-C-N	-5.57	113.01	119.47
1	Q	99	GLN	C-N-CA	-5.57	113.01	119.47
1	P	67	THR	CA-C-N	5.37	127.79	120.54
1	P	67	THR	C-N-CA	5.37	127.79	120.54
1	Q	346	LYS	N-CA-C	-5.14	105.75	111.36
1	Q	285	VAL	N-CA-C	5.09	115.30	108.17

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	P	18	ARG	Sidechain

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	P	4876	0	4819	36	0
1	Q	4859	0	4811	42	0
2	A	32	0	24	12	0
2	B	32	0	24	16	0
3	P	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	Q	2	0	0	0	0
4	P	15	0	0	4	0
4	Q	20	0	0	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	P	167	0	0	0	0
5	Q	183	0	0	3	0
All	All	10190	0	9678	76	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 4.

All (76) 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:Q:346:LYS:HE3	5:Q:938:HOH:O	1.33	1.28
1:P:381:ILE:HD11	2:A:1:01B:H11	1.41	1.01
1:Q:381:ILE:HD11	2:B:1:01B:H11	1.45	0.95
1:P:381:ILE:CD1	2:A:1:01B:H11	1.95	0.95
1:P:275:ALA:O	1:P:279:SER:HB2	1.66	0.94
1:Q:381:ILE:CD1	2:B:1:01B:H11	1.97	0.94
1:Q:378:PHE:HB2	2:B:1:01B:H9	1.63	0.77
1:P:381:ILE:HD11	2:A:1:01B:C11	2.16	0.74
1:Q:381:ILE:HD11	2:B:1:01B:C11	2.21	0.70
1:Q:343:TRP:CE2	1:Q:347:GLU:HG3	2.26	0.70
1:Q:347:GLU:O	1:Q:352:LYS:HB2	1.90	0.69
1:Q:394:LYS:HZ3	2:B:1:01B:H10	1.58	0.69
1:P:378:PHE:HB2	2:A:1:01B:H9	1.75	0.69
1:P:394:LYS:NZ	2:A:1:01B:H10	2.09	0.68
1:Q:378:PHE:CD2	2:B:1:01B:C9	2.81	0.64
1:Q:190:LYS:HG3	4:Q:704:SO4:O1	1.98	0.63
1:P:381:ILE:HD12	2:A:1:01B:H11	1.80	0.62
1:Q:394:LYS:NZ	2:B:1:01B:H10	2.13	0.62
1:Q:381:ILE:HD12	2:B:1:01B:H11	1.79	0.62
1:Q:57:PHE:HA	1:Q:86:GLN:OE1	1.99	0.61
1:P:23:THR:CG2	1:P:146:VAL:HG21	2.31	0.60
1:P:14:PHE:CD2	1:P:28:MET:O	2.54	0.60
1:P:378:PHE:CD2	2:A:1:01B:C9	2.86	0.59
1:Q:484:GLY:H	2:B:4:ALA:HA	1.67	0.58
1:P:23:THR:HG21	1:P:146:VAL:HG21	1.85	0.58
1:Q:28:MET:HE1	1:Q:31:TYR:HB2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:69:ARG:C	1:Q:70:GLU:HG3	2.29	0.56
1:Q:496:HIS:NE2	2:B:1:01B:C6	2.70	0.55
1:P:339:GLU:HG2	1:P:605:GLU:HG2	1.88	0.55
1:P:381:ILE:HB	1:P:413:ASP:HB3	1.89	0.54
1:P:394:LYS:HZ3	2:A:1:01B:H10	1.73	0.53
1:Q:346:LYS:CE	5:Q:938:HOH:O	2.15	0.53
1:Q:343:TRP:CZ2	1:Q:347:GLU:HG3	2.44	0.53
1:Q:381:ILE:HB	1:Q:413:ASP:HB3	1.90	0.53
1:P:394:LYS:NZ	2:A:1:01B:C10	2.71	0.52
1:Q:594:GLY:HA2	1:Q:606:LEU:HD11	1.90	0.52
1:Q:231:ALA:HB3	1:Q:234:GLU:HG2	1.92	0.52
1:Q:279:SER:OG	1:Q:281:GLU:HG3	2.10	0.51
1:Q:394:LYS:NZ	2:B:1:01B:C10	2.74	0.51
1:P:22:LEU:C	1:P:23:THR:CG2	2.84	0.50
1:P:22:LEU:O	1:P:23:THR:HG22	2.11	0.50
1:Q:378:PHE:HB2	2:B:1:01B:C9	2.37	0.49
1:P:367:ARG:HD3	1:P:416:ALA:HB2	1.95	0.49
1:Q:484:GLY:N	2:B:4:ALA:HA	2.28	0.49
1:P:147:SER:C	1:P:148:ILE:HD13	2.38	0.48
1:P:484:GLY:H	2:A:4:ALA:HA	1.78	0.48
1:P:444:LEU:HD11	1:P:477:LEU:HD12	1.96	0.48
1:Q:452:LEU:HD13	1:Q:469:LEU:HD12	1.95	0.48
1:P:140:ALA:HB3	1:Q:284:MET:HE1	1.95	0.47
1:Q:367:ARG:HD3	1:Q:416:ALA:HB2	1.97	0.47
1:P:190:LYS:HG3	4:P:703:SO4:O3	2.15	0.46
1:Q:483:HIS:CG	2:B:2:PRO:HB3	2.51	0.45
1:P:205:ASP:HB3	1:P:494:ASN:ND2	2.32	0.45
1:P:68:ASP:N	1:P:68:ASP:OD1	2.38	0.45
1:Q:231:ALA:HB3	1:Q:234:GLU:CG	2.47	0.45
1:P:310:LEU:O	4:P:703:SO4:O1	2.35	0.44
1:Q:496:HIS:NE2	2:B:1:01B:O	2.50	0.44
1:P:22:LEU:C	1:P:23:THR:HG23	2.42	0.44
1:Q:280:LYS:HE3	1:Q:280:LYS:HB2	1.81	0.44
1:Q:594:GLY:CA	1:Q:606:LEU:HD11	2.48	0.44
1:P:390:LEU:CD2	4:P:705:SO4:O4	2.66	0.43
1:Q:594:GLY:HA2	1:Q:606:LEU:CD1	2.47	0.43
1:P:20:LEU:O	1:P:24:SER:O	2.36	0.43
1:P:465:ARG:O	1:Q:468:THR:OG1	2.32	0.43
1:Q:583:ASN:HD21	1:Q:616:ILE:HG13	1.84	0.43
1:P:78:ARG:CD	2:A:2:PRO:HB2	2.49	0.43
1:P:483:HIS:CG	2:A:2:PRO:HB3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:19:VAL:HG11	1:Q:28:MET:HG3	2.00	0.43
1:Q:26:LYS:HE3	5:Q:853:HOH:O	2.19	0.42
1:P:337:LEU:HD23	1:P:562:VAL:HG21	2.01	0.42
1:P:277:GLU:HG3	1:P:283:HIS:CE1	2.54	0.42
1:P:22:LEU:O	1:P:23:THR:CG2	2.69	0.41
1:Q:470:THR:HG21	1:Q:485:THR:HG23	2.03	0.41
1:Q:378:PHE:CG	2:B:1:01B:C9	3.04	0.41
1:P:85:ASN:HB2	4:P:705:SO4:O1	2.21	0.41
1:Q:24:SER:O	1:Q:25:SER:HB2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	P	611/639 (96%)	589 (96%)	20 (3%)	2 (0%)	36	34
1	Q	608/639 (95%)	590 (97%)	18 (3%)	0	100	100
2	A	2/5 (40%)	0	2 (100%)	0	100	100
2	B	2/5 (40%)	1 (50%)	1 (50%)	0	100	100
All	All	1223/1288 (95%)	1180 (96%)	41 (3%)	2 (0%)	43	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	P	24	SER
1	P	150	GLY

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	P	527/553 (95%)	512 (97%)	15 (3%)	38	40
1	Q	527/553 (95%)	512 (97%)	15 (3%)	38	40
2	A	2/2 (100%)	2 (100%)	0	100	100
2	B	2/2 (100%)	2 (100%)	0	100	100
All	All	1058/1110 (95%)	1028 (97%)	30 (3%)	38	40

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	P	9	LYS
1	P	22	LEU
1	P	24	SER
1	P	68	ASP
1	P	114	GLU
1	P	144	GLN
1	P	270	SER
1	P	279	SER
1	P	280	LYS
1	P	284	MET
1	P	346	LYS
1	P	398	GLU
1	P	425	VAL
1	P	482	GLU
1	P	505	ARG
1	Q	26	LYS
1	Q	68	ASP
1	Q	101	ASP
1	Q	114	GLU
1	Q	173	VAL
1	Q	241	ASN
1	Q	258	SER
1	Q	280	LYS
1	Q	346	LYS

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Mol	Chain	Res	Type
1	Q	394	LYS
1	Q	396	LEU
1	Q	425	VAL
1	Q	528	LYS
1	Q	588	ARG
1	Q	606	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	P	61	ASN
1	P	97	GLN
1	P	283	HIS
1	P	324	GLN
1	P	596	HIS
1	Q	144	GLN
1	Q	324	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	01B	A	1	3,2	11,13,14	1.57	1 (9%)	12,16,18	0.82	0
2	01B	B	1	3,2	11,13,14	1.57	1 (9%)	12,16,18	0.83	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	01B	A	1	3,2	-	1/9/10/12	0/1/1/1
2	01B	B	1	3,2	-	0/9/10/12	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	01B	C6-C7	-5.04	1.39	1.51
2	B	1	01B	C6-C7	-5.03	1.39	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1	01B	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	01B	9	0
2	B	1	01B	13	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	$\# Z > 2$
4	SO4	P	705	-	4,4,4	0.25	0	6,6,6	0.81	0
4	SO4	P	704	-	4,4,4	0.49	0	6,6,6	0.59	0
4	SO4	Q	705	-	4,4,4	0.45	0	6,6,6	0.48	0
4	SO4	Q	706	-	4,4,4	1.29	0	6,6,6	0.93	0
4	SO4	Q	704	-	4,4,4	0.19	0	6,6,6	1.13	0
4	SO4	P	703	-	4,4,4	0.36	0	6,6,6	1.29	0
4	SO4	Q	703	-	4,4,4	1.12	0	6,6,6	1.18	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	P	705	SO4	2	0
4	Q	704	SO4	1	0
4	P	703	SO4	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

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	P	615/639 (96%)	0.29	25 (4%)	41 46	20, 35, 57, 88	5 (0%)
1	Q	612/639 (95%)	0.17	19 (3%)	51 55	15, 32, 53, 87	3 (0%)
2	A	3/5 (60%)	6.65	3 (100%)	0 0	19, 19, 20, 21	3 (100%)
2	B	3/5 (60%)	6.42	3 (100%)	0 0	16, 16, 16, 17	3 (100%)
All	All	1233/1288 (95%)	0.26	50 (4%)	41 46	15, 33, 55, 88	14 (1%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	4	ALA	9.9
2	A	4	ALA	7.5
2	A	2	PRO	6.8
2	B	2	PRO	5.7
2	A	3	PRO	5.6
1	P	508	PRO	4.8
1	Q	1	MET	4.7
1	Q	101	ASP	4.6
1	P	23	THR	4.5
1	Q	509	THR	4.3
1	P	0	HIS	4.2
1	Q	606	LEU	3.9
1	P	25	SER	3.7
1	P	509	THR	3.7
2	B	3	PRO	3.6
1	P	135	PHE	3.4
1	Q	68	ASP	3.3
1	P	616	ILE	3.3
1	P	278	ALA	3.2
1	Q	87	LEU	3.2
1	P	1	MET	3.1

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Mol	Chain	Res	Type	RSRZ
1	P	26	LYS	3.0
1	P	24	SER	2.9
1	Q	275	ALA	2.9
1	P	510	GLY	2.9
1	Q	601	GLY	2.9
1	P	601	GLY	2.8
1	Q	616	ILE	2.7
1	P	21	ALA	2.6
1	P	2	THR	2.6
1	Q	2	THR	2.6
1	Q	135	PHE	2.6
1	Q	278	ALA	2.6
1	Q	510	GLY	2.5
1	Q	4	LEU	2.5
1	Q	69	ARG	2.5
1	P	189	GLU	2.4
1	P	162	ARG	2.4
1	Q	604	ASP	2.4
1	P	280	LYS	2.4
1	Q	607	LYS	2.4
1	P	174	GLU	2.4
1	P	9	LYS	2.3
1	Q	274	LYS	2.3
1	P	163	LEU	2.3
1	P	273	LEU	2.3
1	P	8	ALA	2.3
1	P	279	SER	2.1
1	Q	603	THR	2.1
1	P	173	VAL	2.1

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

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
2	01B	B	1	13/14	0.73	0.43	38,43,53,53	13
2	01B	A	1	13/14	0.77	0.33	39,45,46,46	13

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
4	SO4	P	704	5/5	0.81	0.25	15,22,26,26	0
4	SO4	P	703	5/5	0.93	0.14	17,24,30,31	0
3	ZN	Q	702	1/1	0.97	0.11	48,48,48,48	0
3	ZN	P	701	1/1	0.97	0.12	47,47,47,47	0
3	ZN	P	702	1/1	0.97	0.08	55,55,55,55	0
3	ZN	Q	701	1/1	0.98	0.08	48,48,48,48	0
4	SO4	P	705	5/5	0.98	0.10	20,26,32,32	0
4	SO4	Q	703	5/5	0.98	0.13	12,18,22,24	0
4	SO4	Q	704	5/5	0.98	0.09	23,30,30,34	0
4	SO4	Q	705	5/5	0.98	0.10	17,25,30,31	0
4	SO4	Q	706	5/5	0.98	0.11	10,15,18,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.