Full wwPDB X-ray Structure Validation Report (i)

Oct 7, 2023 – 04:02 PM EDT

PDB ID	:	6DDL
Title	:	Crystal structure of the single mutant (D52N) of NT5C2-Q523X in the basal state
Authors	:	Forouhar, F.; Dieck, C.L.; Tzoneva, G.; Carpenter, Z.; Ambesi-Impiombato, A.; Sanchez-Martin, M.; Kirschner-Schwabe, R.; Lew, S.; Seetharaman, J.; Ferrando, A.A.; Tong, L.
Deposited on	:	2018-05-10
Resolution	:	2.26 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 (i) 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 (i)) were used in the production of this report:

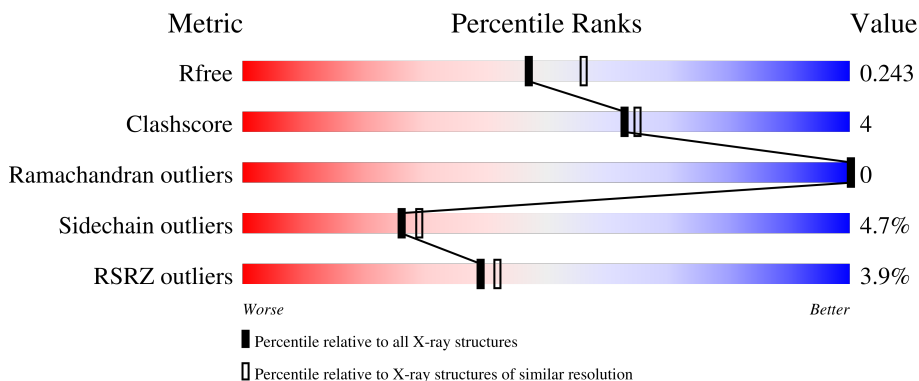|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.26 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	Percentile Ranks	Value
Rfree		0.243
Clashscore		4
Ramachandran outliers		0
Sidechain outliers		4.7%
RSRZ outliers		3.9%

Worse — *Better*

▮ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	541	2% · 74% · 11% · 14%
1	B	541	4% · 74% · 11% · · 13%

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 8292 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 Cytosolic purine 5'-nucleotidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total 3783	C 2444	N 626	O 694	S 19	0	0	0
1	B	469	Total 3835	C 2476	N 640	O 700	S 19	0	0	0

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	GLY	-	expression tag	UNP P49902
A	-16	SER	-	expression tag	UNP P49902
A	-15	SER	-	expression tag	UNP P49902
A	-14	HIS	-	expression tag	UNP P49902
A	-13	HIS	-	expression tag	UNP P49902
A	-12	HIS	-	expression tag	UNP P49902
A	-11	HIS	-	expression tag	UNP P49902
A	-10	HIS	-	expression tag	UNP P49902
A	-9	HIS	-	expression tag	UNP P49902
A	-8	SER	-	expression tag	UNP P49902
A	-7	SER	-	expression tag	UNP P49902
A	-6	GLY	-	expression tag	UNP P49902
A	-5	LEU	-	expression tag	UNP P49902
A	-4	VAL	-	expression tag	UNP P49902
A	-3	PRO	-	expression tag	UNP P49902
A	-2	ARG	-	expression tag	UNP P49902
A	-1	GLY	-	expression tag	UNP P49902
A	0	SER	-	expression tag	UNP P49902
A	52	ASN	ASP	engineered mutation	UNP P49902
B	-17	GLY	-	expression tag	UNP P49902
B	-16	SER	-	expression tag	UNP P49902
B	-15	SER	-	expression tag	UNP P49902
B	-14	HIS	-	expression tag	UNP P49902
B	-13	HIS	-	expression tag	UNP P49902
B	-12	HIS	-	expression tag	UNP P49902

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	HIS	-	expression tag	UNP P49902
B	-10	HIS	-	expression tag	UNP P49902
B	-9	HIS	-	expression tag	UNP P49902
B	-8	SER	-	expression tag	UNP P49902
B	-7	SER	-	expression tag	UNP P49902
B	-6	GLY	-	expression tag	UNP P49902
B	-5	LEU	-	expression tag	UNP P49902
B	-4	VAL	-	expression tag	UNP P49902
B	-3	PRO	-	expression tag	UNP P49902
B	-2	ARG	-	expression tag	UNP P49902
B	-1	GLY	-	expression tag	UNP P49902
B	0	SER	-	expression tag	UNP P49902
B	52	ASN	ASP	engineered mutation	UNP P49902

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).
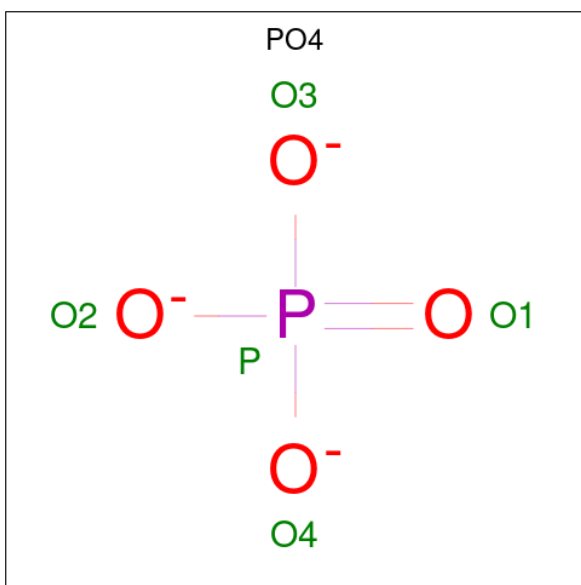
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 5	O 4	P 1	0	0
2	A	1	Total 5	O 4	P 1	0	0
2	B	1	Total 5	O 4	P 1	0	0
2	B	1	Total 5	O 4	P 1	0	0

- Molecule 3 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	336	Total O 336　336		0	0
3	B	318	Total O 318　318		0	0

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: Cytosolic purine 5'-nucleotidase
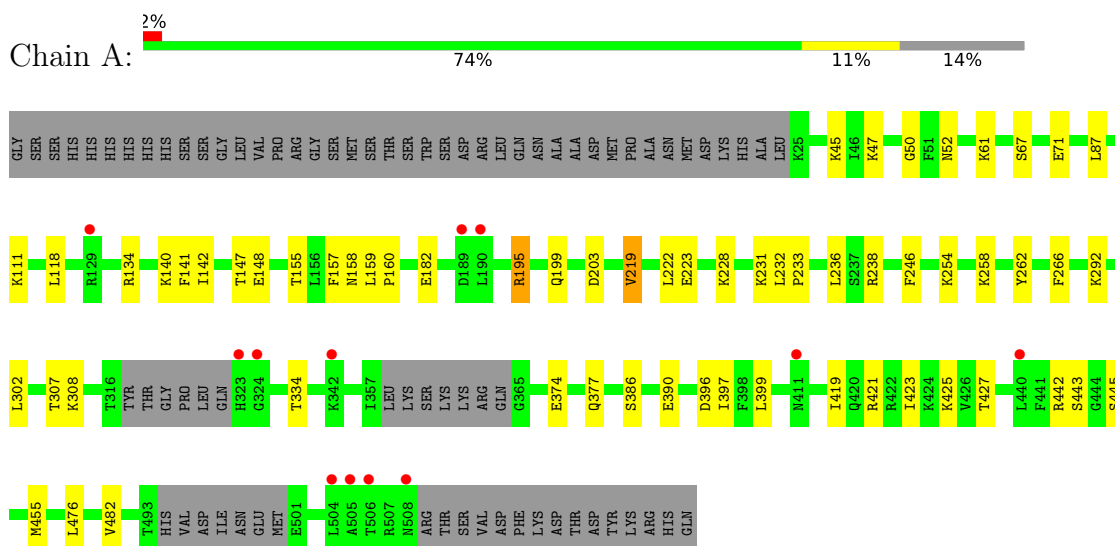
• Molecule 1: Cytosolic purine 5'-nucleotidase
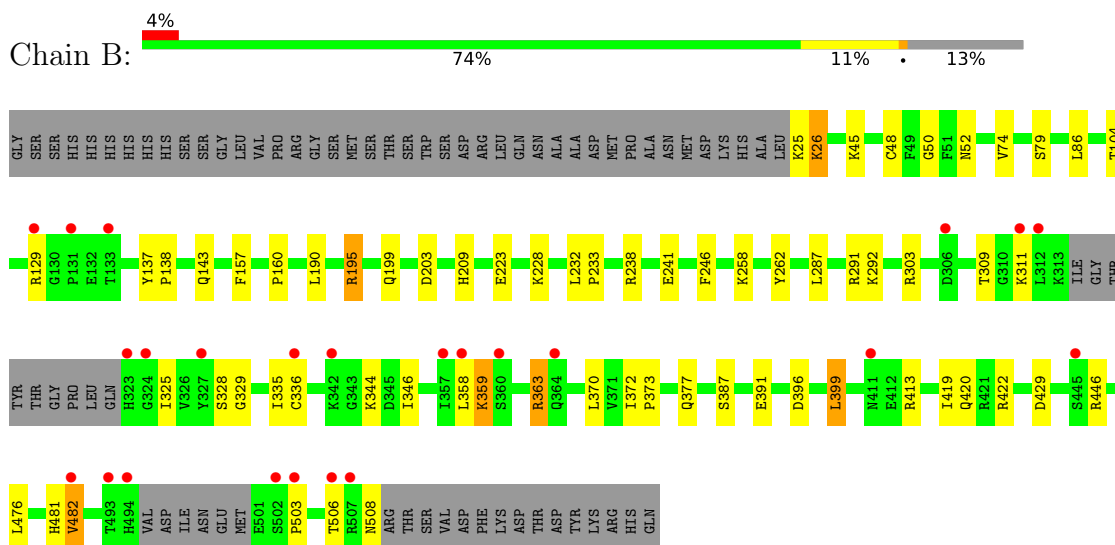
4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	143.45Å 123.63Å 91.25Å 90.00° 115.51° 90.00°	Depositor
Resolution (Å)	49.44 – 2.26 49.44 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.44-2.26) 99.9 (49.44-2.26)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.12 (at 2.27Å)	Xtriage
Refinement program	PHENIX (1.12_2829)	Depositor
R, R_{free}	0.194 , 0.243 0.195 , 0.243	Depositor DCC
R_{free} test set	6782 reflections (10.11%)	wwPDB-VP
Wilson B-factor (Å^2)	34.3	Xtriage
Anisotropy	0.523	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 52.0	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8292	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.29% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: PO4

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	0.38	0/3879	0.56	0/5238
1	B	0.39	0/3933	0.55	0/5308
All	All	0.39	0/7812	0.55	0/10546

There are no bond length outliers.

There are no bond angle outliers.

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	3783	0	3714	34	0
1	B	3835	0	3777	37	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	336	0	0	8	1
3	B	318	0	0	10	1
All	All	8292	0	7491	65	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 4.

All (65) 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:397:ILE:HG12	1:B:26:LYS:HB3	1.72	0.72
1:B:25:LYS:N	3:B:705:HOH:O	2.22	0.71
1:B:287:LEU:HD13	1:B:335:ILE:HD11	1.73	0.70
1:B:429:ASP:OD1	3:B:701:HOH:O	2.12	0.68
1:B:228:LYS:NZ	3:B:707:HOH:O	2.26	0.68
1:A:118:LEU:HD11	1:A:142:ILE:HG23	1.77	0.66
1:B:325:ILE:O	3:B:702:HOH:O	2.13	0.66
1:A:141:PHE:CD1	1:B:359:LYS:HE3	2.32	0.65
1:B:241:GLU:OE2	3:B:703:HOH:O	2.15	0.64
1:A:425:LYS:NZ	3:A:707:HOH:O	2.32	0.62
1:B:195:ARG:NH2	1:B:199:GLN:OE1	2.33	0.61
1:A:421:ARG:NH1	3:A:711:HOH:O	2.35	0.59
1:B:303:ARG:HB2	1:B:328:SER:HB3	1.84	0.59
1:A:302:LEU:HD22	1:A:334:THR:HG21	1.86	0.57
1:A:195:ARG:NH2	1:A:199:GLN:OE1	2.38	0.57
1:B:258:LYS:NZ	3:B:710:HOH:O	2.29	0.57
1:A:67:SER:O	1:A:71:GLU:HG3	2.06	0.56
1:A:482:VAL:HG12	1:B:482:VAL:HG12	1.88	0.55
1:B:373:PRO:HG3	3:B:885:HOH:O	2.06	0.55
1:B:228:LYS:HE3	1:B:262:TYR:O	2.08	0.54
1:A:71:GLU:HG2	3:A:960:HOH:O	2.08	0.53
1:A:52:ASN:OD1	1:A:292:LYS:NZ	2.41	0.53
1:B:203:ASP:OD2	3:B:704:HOH:O	2.18	0.53
1:B:52:ASN:OD1	1:B:292:LYS:NZ	2.41	0.53
1:B:391:GLU:OE2	1:B:422:ARG:NH2	2.43	0.52
1:A:232:LEU:HB3	1:A:233:PRO:HD3	1.92	0.51
1:A:442:ARG:NH2	1:A:445:SER:O	2.43	0.51
1:B:232:LEU:HB3	1:B:233:PRO:HD3	1.91	0.51
1:B:157:PHE:O	1:B:160:PRO:HD2	2.11	0.50
1:A:399:LEU:HD13	1:A:419:ILE:HG21	1.93	0.50
1:B:413:ARG:NH2	3:B:726:HOH:O	2.46	0.49
1:A:219:VAL:HG11	1:A:258:LYS:HD3	1.95	0.48
1:A:134:ARG:O	1:B:363:ARG:NH2	2.46	0.48
1:A:199:GLN:NE2	3:A:717:HOH:O	2.40	0.48
1:A:423:ILE:O	1:A:427:THR:HG23	2.12	0.47
1:A:231:LYS:NZ	3:A:712:HOH:O	2.47	0.47
1:B:287:LEU:HD11	1:B:329:GLY:O	2.14	0.47
1:A:254:LYS:HE3	1:A:254:LYS:HB2	1.67	0.47
1:A:111:LYS:HE2	1:A:443:SER:HB2	1.96	0.47
1:A:134:ARG:HH22	1:B:344:LYS:NZ	2.13	0.46
1:B:129:ARG:HD3	1:B:129:ARG:HA	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:THR:O	1:A:158:ASN:HB2	2.15	0.46
1:B:48:CYS:HB2	1:B:346:ILE:HG12	1.98	0.46
1:A:374:GLU:O	1:A:377:GLN:HG2	2.15	0.45
1:B:50:GLY:HA3	1:B:246:PHE:CZ	2.51	0.45
1:A:140:LYS:NZ	3:A:721:HOH:O	2.47	0.45
1:A:157:PHE:O	1:A:160:PRO:HD2	2.17	0.45
1:B:399:LEU:HD13	1:B:419:ILE:HG21	1.98	0.45
1:A:50:GLY:HA3	1:A:246:PHE:CZ	2.51	0.45
1:B:363:ARG:HE	1:B:363:ARG:HB3	1.64	0.44
1:B:74:VAL:HG13	1:B:86:LEU:HB3	1.98	0.44
1:A:386:SER:O	1:A:390:GLU:HG3	2.18	0.44
1:A:396:ASP:OD1	1:B:26:LYS:NZ	2.44	0.44
1:B:370:LEU:HD23	1:B:372:ILE:HD11	2.00	0.43
1:A:228:LYS:HG2	1:A:266:PHE:CZ	2.53	0.43
1:A:203:ASP:OD2	3:A:701:HOH:O	2.21	0.43
1:B:209:HIS:HE1	3:B:922:HOH:O	2.00	0.43
1:A:61:LYS:NZ	1:A:223:GLU:O	2.37	0.43
1:A:47:LYS:HE3	3:A:957:HOH:O	2.19	0.43
1:B:503:PRO:HA	1:B:506:THR:OG1	2.20	0.42
1:B:363:ARG:H	1:B:363:ARG:HG2	1.58	0.42
1:B:104:THR:HB	1:B:195:ARG:HA	2.02	0.42
1:A:222:LEU:HD13	1:A:262:TYR:HB2	2.03	0.41
1:B:399:LEU:HD12	1:B:399:LEU:HA	1.97	0.41
1:B:137:TYR:HA	1:B:138:PRO:HD2	1.94	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)
3:A:942:HOH:O	3:B:986:HOH:O[4_555]	2.17	0.03

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	456/541 (84%)	448 (98%)	8 (2%)	0	100	100
1	B	463/541 (86%)	452 (98%)	11 (2%)	0	100	100
All	All	919/1082 (85%)	900 (98%)	19 (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	414/483 (86%)	400 (97%)	14 (3%)	37	45
1	B	420/483 (87%)	395 (94%)	25 (6%)	19	18
All	All	834/966 (86%)	795 (95%)	39 (5%)	26	29

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LYS
1	A	87	LEU
1	A	147	THR
1	A	148	GLU
1	A	159	LEU
1	A	182	GLU
1	A	195	ARG
1	A	219	VAL
1	A	236	LEU
1	A	238	ARG
1	A	307	THR
1	A	308	LYS
1	A	455	MET
1	A	476	LEU
1	B	26	LYS
1	B	45	LYS

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Mol	Chain	Res	Type
1	B	79	SER
1	B	143	GLN
1	B	190	LEU
1	B	195	ARG
1	B	223	GLU
1	B	238	ARG
1	B	291	ARG
1	B	309	THR
1	B	311	LYS
1	B	336	CYS
1	B	358	LEU
1	B	359	LYS
1	B	363	ARG
1	B	377	GLN
1	B	387	SER
1	B	396	ASP
1	B	399	LEU
1	B	420	GLN
1	B	446	ARG
1	B	476	LEU
1	B	481	HIS
1	B	482	VAL
1	B	508	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3   RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5   Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 ligands 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	PO4	B	601	-	4,4,4	0.62	0	6,6,6	0.45	0
2	PO4	A	601	-	4,4,4	0.73	0	6,6,6	0.74	0
2	PO4	A	602	-	4,4,4	1.02	0	6,6,6	0.51	0
2	PO4	B	602	-	4,4,4	0.85	0	6,6,6	0.90	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.

No monomer is involved in short contacts.

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 ⓘ

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	464/541 (85%)	0.13	12 (2%)	56	59	23, 36, 66, 98	0
1	B	469/541 (86%)	0.21	24 (5%)	28	30	24, 39, 73, 109	0
All	All	933/1082 (86%)	0.17	36 (3%)	39	42	23, 37, 70, 109	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	129	ARG	4.5
1	B	323	HIS	3.9
1	B	357	ILE	3.8
1	B	324	GLY	3.7
1	B	364	GLN	3.7
1	A	189	ASP	3.7
1	B	133	THR	3.3
1	B	342	LYS	3.1
1	B	358	LEU	3.1
1	A	323	HIS	3.0
1	B	494	HIS	2.9
1	A	190	LEU	2.9
1	A	504	LEU	2.9
1	A	129	ARG	2.8
1	B	506	THR	2.8
1	A	411	ASN	2.7
1	B	503	PRO	2.7
1	B	131	PRO	2.4
1	A	342	LYS	2.4
1	B	311	LYS	2.3
1	B	482	VAL	2.3
1	A	324	GLY	2.3
1	B	445	SER	2.3
1	B	411	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	506	THR	2.2
1	A	505	ALA	2.2
1	B	306	ASP	2.2
1	B	493	THR	2.1
1	B	360	SER	2.1
1	B	312	LEU	2.1
1	B	502	SER	2.1
1	B	327	TYR	2.1
1	A	508	ASN	2.0
1	A	440	LEU	2.0
1	B	507	ARG	2.0
1	B	336	CYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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	PO4	B	601	5/5	0.89	0.19	37,41,52,64	0
2	PO4	A	601	5/5	0.93	0.17	41,47,50,56	0
2	PO4	A	602	5/5	0.97	0.13	29,32,33,39	0
2	PO4	B	602	5/5	0.98	0.10	27,30,34,34	0

6.5 Other polymers ⓘ

There are no such residues in this entry.