



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:19 AM UTC

PDB ID : 3H64 / pdb_00003h64
Title : Catalytic domain of human Serine/Threonine Phosphatase 5 (PP5c) with two Mn²⁺ atoms complexed with endothall
Authors : Bertini, I.; Calderone, V.; Fragai, M.; Luchinat, C.; Talluri, E.
Deposited on : 2009-04-23
Resolution : 1.90 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

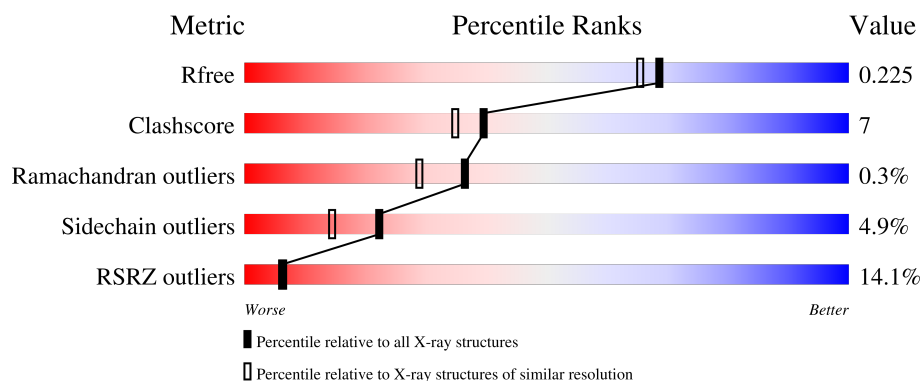
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	315	86% 11% ..
1	D	315	26% 77% 19% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5259 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.

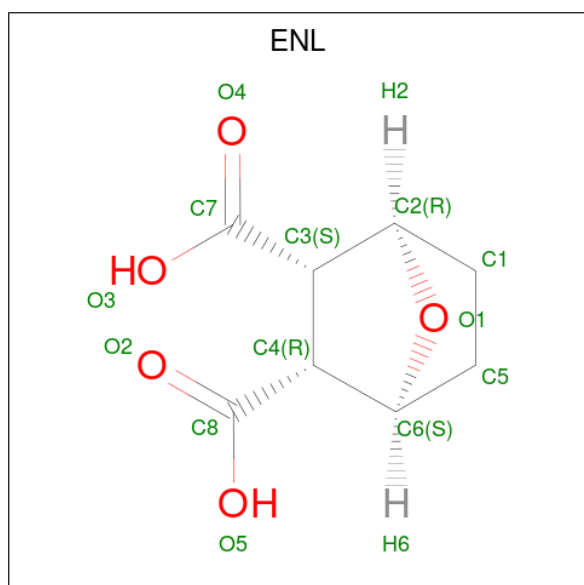
- Molecule 1 is a protein called Serine/threonine-protein phosphatase 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	0	0
			2482	1584	418	466	14			
1	D	309	Total	C	N	O	S	0	0	0
			2482	1584	418	465	15			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		
2	D	2	Total	Mn	0	0
			2	2		

- Molecule 3 is (1R,2S,3R,4S)-7-oxabicyclo[2.2.1]heptane-2,3-dicarboxylic acid (CCD ID: ENL) (formula: C₈H₁₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			26	16	10		
3	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	213	Total	O	0	0
			213	213		
4	D	39	Total	O	0	0
			39	39		

- Molecule 1: Serine/threonine-protein phosphatase 5



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.55Å 41.46Å 95.61Å 90.00° 96.78° 90.00°	Depositor
Resolution (Å)	38.38 – 1.90 38.38 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (38.38-1.90) 99.9 (38.38-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.62 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.5.0089	Depositor
R, R_{free}	0.190 , 0.237 (Not available) , 0.225	Depositor DCC
R_{free} test set	4389 reflections (9.14%)	wwPDB-VP
Wilson B-factor (Å ²)	16.6	Xtriage
Anisotropy	0.688	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5259	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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: MN, ENL

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	1.09	1/2544 (0.0%)	1.03	2/3441 (0.1%)
1	D	0.69	0/2544	0.97	6/3441 (0.2%)
All	All	0.91	1/5088 (0.0%)	1.00	8/6882 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	479	VAL	CA-CB	6.02	1.60	1.55

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	PRO	CA-C-N	-6.21	114.33	120.98
1	A	375	PRO	C-N-CA	-6.21	114.33	120.98
1	D	427	HIS	N-CA-CB	-5.75	107.62	114.17
1	D	407	GLY	CA-C-N	5.59	126.83	119.84
1	D	407	GLY	C-N-CA	5.59	126.83	119.84
1	D	314	GLY	N-CA-C	5.31	119.26	113.58
1	D	244	HIS	N-CA-C	5.19	118.40	111.24
1	D	247	PHE	N-CA-C	5.08	116.50	111.07

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	2482	0	2414	26	0
1	D	2482	0	2414	39	0
2	A	2	0	0	0	0
2	D	2	0	0	0	0
3	A	26	0	16	5	0
3	D	13	0	8	5	0
4	A	213	0	0	8	0
4	D	39	0	0	2	0
All	All	5259	0	4852	66	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 7.

All (66) 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:D:275:ARG:HE	3:D:0:ENL:H5	1.23	1.03
1:D:188:ILE:CD1	1:D:192:LYS:HE3	1.91	1.00
1:D:275:ARG:NE	3:D:0:ENL:H5	1.84	0.93
3:A:0[B]:ENL:H2	4:A:60:HOH:O	1.71	0.90
1:D:188:ILE:HD11	1:D:192:LYS:HE3	1.60	0.82
1:D:188:ILE:HD12	1:D:192:LYS:HE3	1.64	0.79
1:D:415:LEU:HD11	1:D:441:ARG:HB3	1.67	0.75
1:A:474:HIS:HE1	4:A:549:HOH:O	1.71	0.74
1:D:188:ILE:HB	4:D:152:HOH:O	1.92	0.70
1:D:419:ASN:N	1:D:419:ASN:HD22	1.93	0.66
1:A:326:GLN:HG3	4:A:79:HOH:O	1.96	0.65
1:A:465:GLN:HB2	4:A:506:HOH:O	1.95	0.65
1:A:205:HIS:HD2	1:A:207:LYS:H	1.45	0.62
1:D:307:ASP:O	1:D:311:GLN:HG2	2.00	0.61
1:A:205:HIS:CD2	1:A:207:LYS:HG3	2.36	0.60
1:A:298:HIS:HE1	4:A:49:HOH:O	1.84	0.60
1:D:222:LYS:HB2	4:D:511:HOH:O	2.00	0.59
1:D:415:LEU:CD1	1:D:441:ARG:HB3	2.31	0.58
1:A:451:TYR:OH	3:A:0[B]:ENL:H5A	2.03	0.58
1:D:436:VAL:HG12	1:D:440:GLY:HA2	1.85	0.57
1:D:309:MET:HE3	1:D:386:TRP:HZ2	1.71	0.56
1:D:487:MET:C	1:D:489:TYR:H	2.14	0.56
1:D:309:MET:HE3	1:D:386:TRP:CZ2	2.41	0.55
1:A:217:LYS:HB2	1:A:334:VAL:HG22	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:304:HIS:NE2	3:D:0:ENL:H1	2.23	0.53
1:D:330:LEU:O	1:D:334:VAL:HG23	2.08	0.53
1:A:205:HIS:CD2	1:A:207:LYS:H	2.24	0.53
1:A:388:ASP:O	1:A:405:GLN:HA	2.09	0.51
1:A:273:VAL:O	1:A:274:ASP:HB2	2.12	0.49
1:A:474:HIS:CE1	4:A:549:HOH:O	2.53	0.49
1:D:427:HIS:CE1	3:D:0:ENL:O3	2.66	0.48
1:D:332:SER:O	1:D:336:GLU:HG3	2.13	0.48
1:A:298:HIS:HD2	4:A:38:HOH:O	1.97	0.48
1:A:447:SER:HA	1:A:459:ALA:HB1	1.95	0.48
1:A:329:GLU:HB3	4:A:167:HOH:O	2.14	0.47
1:A:275:ARG:NH1	1:A:304:HIS:CE1	2.83	0.47
1:D:267:ILE:HG12	1:D:298:HIS:HB2	1.96	0.47
1:D:451:TYR:OH	3:D:0:ENL:H6	2.13	0.47
1:A:275:ARG:HG3	1:A:451:TYR:OH	2.13	0.47
1:A:329:GLU:O	1:A:333:GLU:HG3	2.15	0.47
1:D:388:ASP:O	1:D:405:GLN:HA	2.16	0.46
1:A:399:LYS:HD3	1:A:405:GLN:OE1	2.14	0.46
1:D:417:GLU:O	1:D:417:GLU:HG2	2.15	0.46
1:A:275:ARG:CZ	3:A:0[B]:ENL:H5	2.46	0.46
1:A:451:TYR:OH	3:A:0[B]:ENL:C5	2.64	0.46
1:D:263:THR:C	1:D:265:PRO:HD3	2.40	0.46
1:D:390:GLN:HB3	1:D:405:GLN:HE21	1.82	0.45
1:D:237:ILE:HG13	1:D:265:PRO:HB2	1.99	0.44
1:D:225:THR:HG21	1:D:369:ILE:HB	1.99	0.44
1:A:382:CYS:SG	1:A:402:VAL:HG23	2.58	0.44
1:D:449:PRO:O	1:D:450:ASN:C	2.61	0.43
1:D:273:VAL:O	1:D:274:ASP:HB2	2.18	0.43
1:D:244:HIS:HE1	1:D:304:HIS:CD2	2.37	0.42
1:A:244:HIS:O	1:A:276:GLY:HA3	2.19	0.42
1:D:249:ASP:O	1:D:253:ILE:HG13	2.18	0.42
1:D:246:GLN:HE22	1:D:451:TYR:HA	1.83	0.41
1:D:399:LYS:HE2	1:D:399:LYS:HB3	1.75	0.41
1:A:248:TYR:CE1	1:A:482:PRO:HD2	2.55	0.41
1:A:246:GLN:HE22	1:A:451:TYR:HA	1.84	0.41
1:D:232:LYS:HE2	1:D:232:LYS:HB3	1.58	0.41
1:A:427:HIS:CE1	3:A:0[B]:ENL:O3	2.74	0.41
1:D:267:ILE:HA	1:D:298:HIS:O	2.21	0.41
1:D:403:SER:OG	1:D:404:CYS:N	2.53	0.41
1:D:182:GLU:O	1:D:183:ASP:HB2	2.20	0.40
1:D:242:ASP:HB3	1:D:244:HIS:CD2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:279:SER:HB2	1:D:318:GLU:OE1	2.22	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	A	307/315 (98%)	297 (97%)	10 (3%)	0	100	100
1	D	307/315 (98%)	278 (91%)	27 (9%)	2 (1%)	18	10
All	All	614/630 (98%)	575 (94%)	37 (6%)	2 (0%)	36	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	308	ASN
1	D	488	ALA

5.3.2 Protein sidechains [i](#)

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	275/279 (99%)	265 (96%)	10 (4%)	31	23
1	D	274/279 (98%)	257 (94%)	17 (6%)	16	9
All	All	549/558 (98%)	522 (95%)	27 (5%)	22	14

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	188	ILE
1	A	216	VAL
1	A	231	LEU
1	A	233	GLU
1	A	399	LYS
1	A	402	VAL
1	A	419	ASN
1	A	425	ARG
1	A	465	GLN
1	A	484	VAL
1	D	188	ILE
1	D	189	SER
1	D	232	LYS
1	D	233	GLU
1	D	255	GLU
1	D	262	GLU
1	D	311	GLN
1	D	364	ASP
1	D	370	GLU
1	D	377	ASP
1	D	397	ILE
1	D	399	LYS
1	D	402	VAL
1	D	419	ASN
1	D	425	ARG
1	D	429	VAL
1	D	484	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	201	GLN
1	A	205	HIS
1	A	246	GLN
1	A	264	ASN
1	A	298	HIS
1	A	310	ASN
1	A	311	GLN
1	A	393	ASN
1	A	465	GLN
1	D	205	HIS

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Mol	Chain	Res	Type
1	D	246	GLN
1	D	264	ASN
1	D	298	HIS
1	D	310	ASN
1	D	311	GLN
1	D	326	GLN
1	D	405	GLN
1	D	419	ASN
1	D	475	GLN

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](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\# Z > 2$
3	ENL	A	0[B]	2	14,14,14	1.39	1 (7%)	21,21,21	1.87	5 (23%)
3	ENL	A	0[A]	2	14,14,14	1.01	0	21,21,21	1.76	3 (14%)
3	ENL	D	0	2	14,14,14	1.13	1 (7%)	21,21,21	1.63	4 (19%)

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
3	ENL	A	0[B]	2	-	0/8/29/29	0/3/2/2
3	ENL	A	0[A]	2	-	0/8/29/29	0/3/2/2
3	ENL	D	0	2	-	0/8/29/29	0/3/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	0[B]	ENL	C3-C2	-2.58	1.50	1.54
3	D	0	ENL	O3-C7	-2.08	1.24	1.30

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	0[A]	ENL	C1-C2-C3	-5.25	104.07	109.69
3	A	0[B]	ENL	C2-C3-C7	-5.20	96.54	111.96
3	D	0	ENL	C4-C3-C7	4.22	124.80	115.07
3	A	0[A]	ENL	C3-C4-C8	-3.63	106.69	115.07
3	D	0	ENL	C2-C3-C7	-3.47	101.66	111.96
3	A	0[B]	ENL	O2-C8-C4	-3.13	114.86	122.84
3	A	0[B]	ENL	O5-C8-C4	2.42	120.86	113.91
3	A	0[B]	ENL	O1-C6-C5	-2.40	100.28	104.34
3	A	0[B]	ENL	O1-C2-C3	-2.39	98.55	103.34
3	A	0[A]	ENL	O3-C7-C3	2.21	120.28	113.91
3	D	0	ENL	O1-C6-C5	-2.09	100.81	104.34
3	D	0	ENL	O2-C8-C4	-2.02	117.67	122.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	0[B]	ENL	5	0
3	D	0	ENL	5	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	A	309/315 (98%)	-0.51	4 (1%) 75 78	3, 12, 26, 50	2 (0%)
1	D	309/315 (98%)	1.45	83 (26%) 1 1	21, 43, 61, 72	3 (0%)
All	All	618/630 (98%)	0.47	87 (14%) 6 6	3, 27, 58, 72	5 (0%)

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	176	TYR	7.1
1	A	484	VAL	5.1
1	D	231	LEU	4.9
1	D	312	ILE	4.4
1	D	208	CYS	3.9
1	D	416	GLU	3.8
1	D	490	ALA	3.7
1	D	479	VAL	3.7
1	D	473	PHE	3.7
1	D	319	VAL	3.6
1	D	413	ALA	3.5
1	D	360	GLY	3.4
1	D	348	VAL	3.3
1	D	484	VAL	3.3
1	D	482	PRO	3.2
1	D	476	PHE	3.2
1	D	462	ILE	3.1
1	D	210	TYR	3.0
1	D	488	ALA	3.0
1	D	469	LEU	3.0
1	D	346	GLY	3.0
1	D	315	PHE	2.9
1	D	466	GLY	2.9
1	D	197	TRP	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	455	MET	2.9
1	D	313	TYR	2.8
1	D	184	GLY	2.8
1	D	204	LEU	2.8
1	D	188	ILE	2.8
1	D	361	VAL	2.8
1	D	182	GLU	2.7
1	A	231	LEU	2.7
1	D	478	ALA	2.7
1	D	377	ASP	2.7
1	D	464	LEU	2.7
1	D	232	LYS	2.7
1	D	237	ILE	2.6
1	D	487	MET	2.6
1	D	461	TYR	2.6
1	D	250	LEU	2.6
1	D	311	GLN	2.6
1	D	209	ALA	2.6
1	D	187	THR	2.6
1	D	276	GLY	2.6
1	D	474	HIS	2.5
1	D	247	PHE	2.5
1	D	363	LEU	2.5
1	D	330	LEU	2.5
1	D	411	THR	2.5
1	D	489	TYR	2.4
1	D	477	THR	2.4
1	D	480	PRO	2.4
1	D	186	VAL	2.4
1	D	397	ILE	2.4
1	D	452	CYS	2.4
1	D	230	THR	2.3
1	D	194	LEU	2.3
1	D	234	THR	2.3
1	D	459	ALA	2.3
1	D	321	ALA	2.3
1	D	471	PRO	2.3
1	D	433	GLY	2.3
1	D	481	HIS	2.3
1	D	317	GLY	2.2
1	D	467	SER	2.2
1	D	258	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	265	PRO	2.2
1	D	198	TYR	2.2
1	D	189	SER	2.2
1	A	177	SER	2.2
1	D	463	HIS	2.1
1	D	254	PHE	2.1
1	D	414	PHE	2.1
1	D	314	GLY	2.1
1	D	486	PRO	2.1
1	D	278	PHE	2.1
1	D	472	GLN	2.1
1	D	456	GLY	2.1
1	D	325	ALA	2.1
1	D	324	THR	2.1
1	D	248	TYR	2.0
1	D	392	GLN	2.0
1	D	263	THR	2.0
1	D	454	GLN	2.0
1	D	308	ASN	2.0
1	D	419	ASN	2.0
1	D	451	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](#)

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
3	ENL	D	0	13/13	0.83	0.15	32,42,45,47	0
3	ENL	A	0[B]	13/13	0.95	0.06	9,12,14,15	13

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ENL	A	0[A]	13/13	0.95	0.06	9,11,15,15	13
2	MN	D	501	1/1	0.99	0.04	10,10,10,10	0
2	MN	D	500	1/1	0.99	0.02	9,9,9,9	0
2	MN	A	501	1/1	1.00	0.01	2,2,2,2	0
2	MN	A	500	1/1	1.00	0.01	2,2,2,2	0

6.5 Other polymers [i](#)

There are no such residues in this entry.