



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 04:16 am BST

PDB ID : 6OBN
Title : The crystal structure of coexpressed SDS22:PP1 complex
Authors : Choy, M.S.; Moon, T.M.; Bray, J.A.; Archuleta, T.L.; Shi, W.; Peti, W.; Page, R.
Deposited on : 2019-03-21
Resolution : 2.70 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

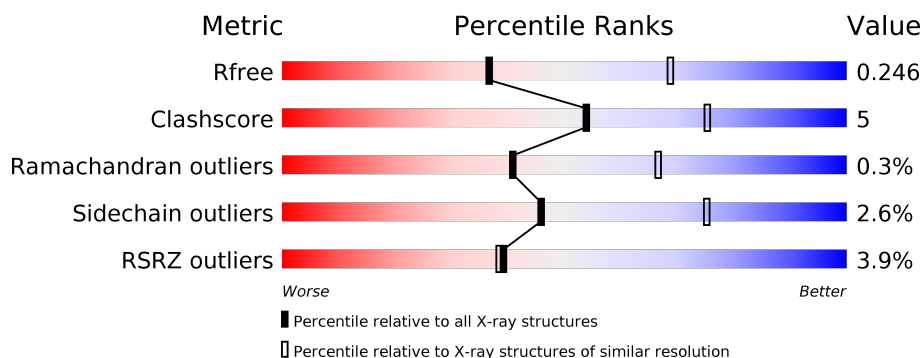
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 on 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	305	<div> <div>2%</div> <div>88%</div> <div>9%</div> <div>••</div> </div>
1	B	305	<div> <div>3%</div> <div>82%</div> <div>13%</div> <div>• 5%</div> </div>
2	C	310	<div> <div>4%</div> <div>78%</div> <div>13%</div> <div>• 8%</div> </div>
2	D	310	<div> <div>5%</div> <div>74%</div> <div>17%</div> <div>• 9%</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PO4	B	401	-	-	X	-
3	PO4	C	403	-	-	-	X
3	PO4	D	401	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9480 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 PP1-alpha catalytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	0	0
			2397	1533	401	444	19			
1	B	291	Total	C	N	O	S	0	0	0
			2332	1495	390	429	18			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P62136
A	-3	HIS	-	expression tag	UNP P62136
A	-2	MET	-	expression tag	UNP P62136
A	-1	GLY	-	expression tag	UNP P62136
A	0	SER	-	expression tag	UNP P62136
B	-4	GLY	-	expression tag	UNP P62136
B	-3	HIS	-	expression tag	UNP P62136
B	-2	MET	-	expression tag	UNP P62136
B	-1	GLY	-	expression tag	UNP P62136
B	0	SER	-	expression tag	UNP P62136

- Molecule 2 is a protein called Protein phosphatase 1 regulatory subunit 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	285	Total	C	N	O	S	0	0	0
			2322	1465	405	445	7			
2	D	283	Total	C	N	O	S	0	0	0
			2300	1451	400	442	7			

There are 10 discrepancies between the modelled and reference sequences:

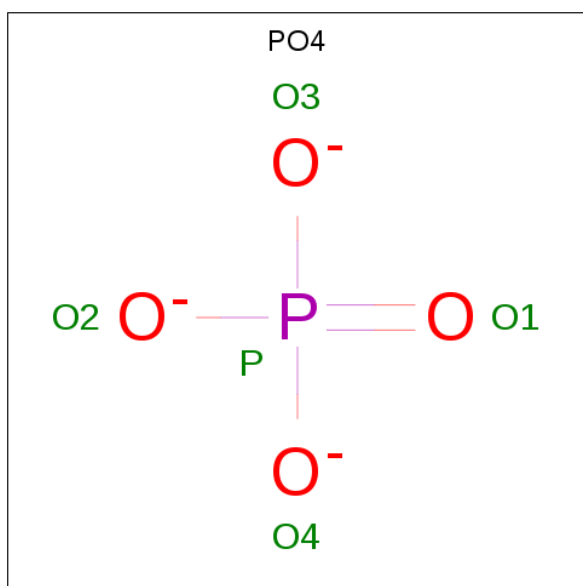
Chain	Residue	Modelled	Actual	Comment	Reference
C	51	GLY	-	expression tag	UNP Q15435

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Chain	Residue	Modelled	Actual	Comment	Reference
C	52	HIS	-	expression tag	UNP Q15435
C	53	MET	-	expression tag	UNP Q15435
C	54	GLY	-	expression tag	UNP Q15435
C	55	SER	-	expression tag	UNP Q15435
D	51	GLY	-	expression tag	UNP Q15435
D	52	HIS	-	expression tag	UNP Q15435
D	53	MET	-	expression tag	UNP Q15435
D	54	GLY	-	expression tag	UNP Q15435
D	55	SER	-	expression tag	UNP Q15435

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	O	P	0	0
			5	4	1		

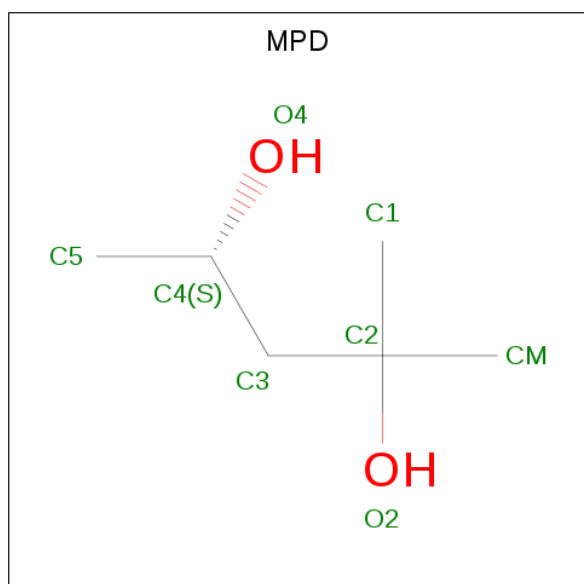
- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Fe	0	0
			1	1		
4	A	1	Total	Fe	0	0
			1	1		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Cl	0	0
			1	1		
5	A	2	Total	Cl	0	0
			2	2		

- Molecule 6 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			8	6	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			8	6	2		
6	C	1	Total	C	O	0	0
			8	6	2		
6	C	1	Total	C	O	0	0
			8	6	2		

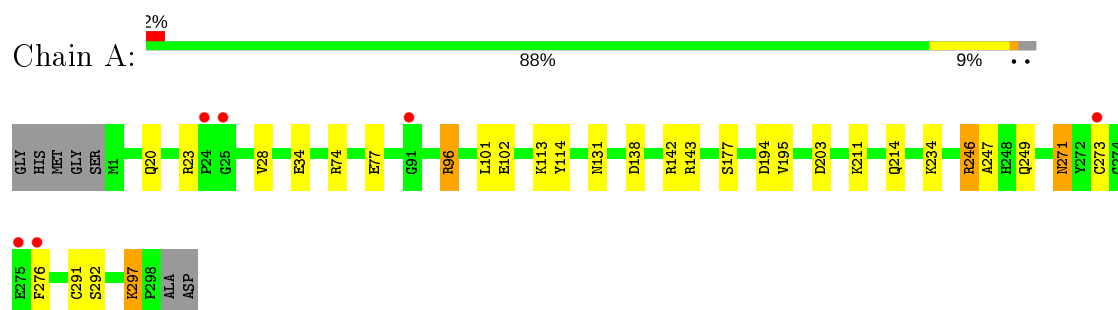
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	20	Total	O	0	0
			20	20		
7	B	12	Total	O	0	0
			12	12		
7	C	13	Total	O	0	0
			13	13		
7	D	7	Total	O	0	0
			7	7		

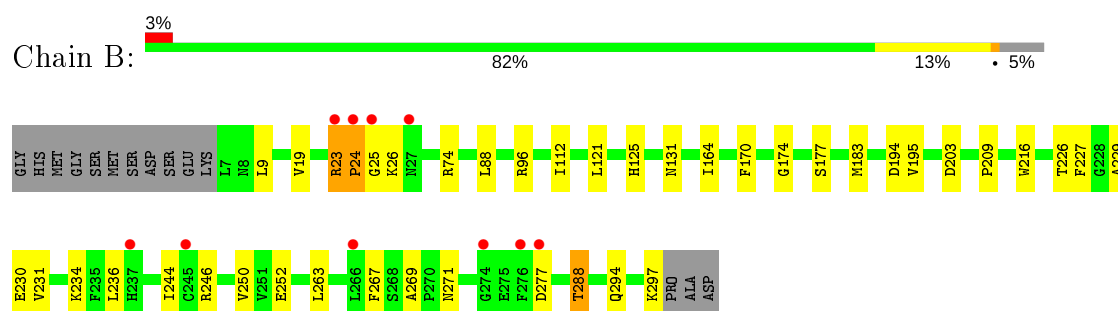
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

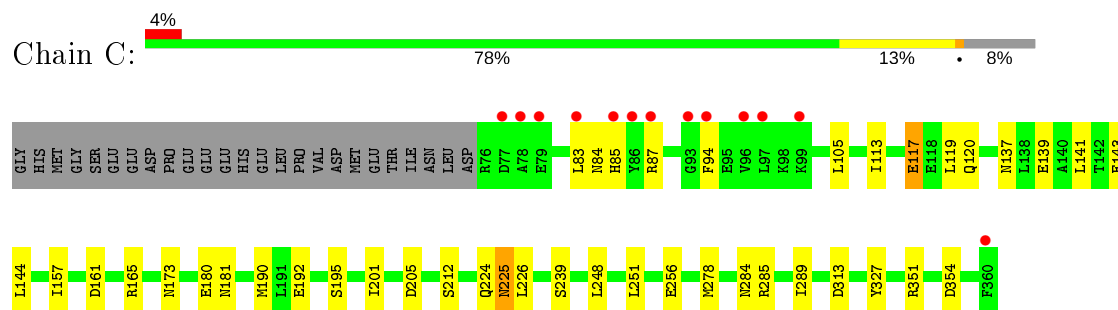
- Molecule 1: Serine/threonine-protein phosphatase PP1-alpha catalytic subunit



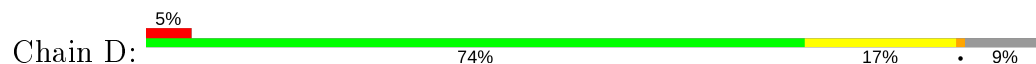
- Molecule 1: Serine/threonine-protein phosphatase PP1-alpha catalytic subunit

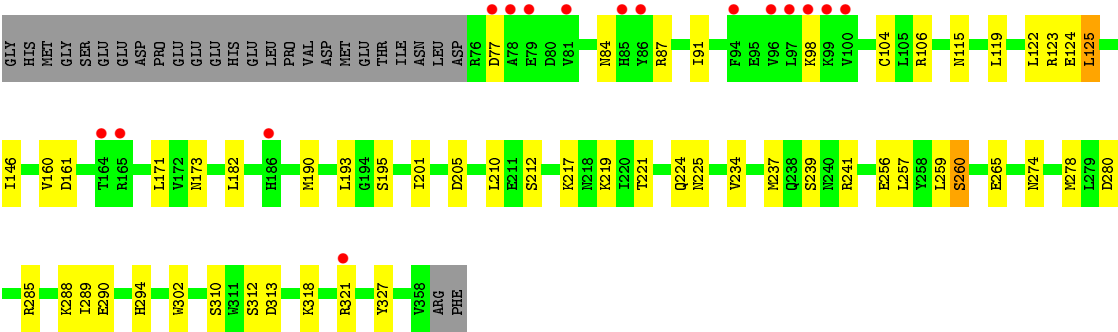


- Molecule 2: Protein phosphatase 1 regulatory subunit 7



- Molecule 2: Protein phosphatase 1 regulatory subunit 7





4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	94.15Å 94.15Å 545.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.87 – 2.70 29.87 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.87-2.70) 92.3 (29.87-2.70)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 2.68Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.202 , 0.246 0.202 , 0.246	Depositor DCC
R_{free} test set	1986 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	52.2	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9480	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MPD, CL, PO4, FE

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	0.32	0/2451	0.50	0/3309
1	B	0.31	0/2385	0.50	0/3223
2	C	0.28	0/2344	0.53	0/3158
2	D	0.28	0/2321	0.54	1/3127 (0.0%)
All	All	0.30	0/9501	0.52	1/12817 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	125	LEU	CA-CB-CG	5.89	128.84	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2397	0	2358	18	0
1	B	2332	0	2283	25	0
2	C	2322	0	2402	24	0
2	D	2300	0	2385	34	0
3	A	10	0	0	0	0
3	B	5	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	15	0	0	0	0
3	D	10	0	0	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
6	A	16	0	28	3	0
6	C	16	0	28	2	0
7	A	20	0	0	1	0
7	B	12	0	0	1	0
7	C	13	0	0	0	0
7	D	7	0	0	0	0
All	All	9480	0	9484	101	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 5.

All (101) 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:B:271:ASN:OD1	1:B:297:LYS:NZ	2.16	0.78
2:D:318:LYS:O	2:D:321:ARG:NE	2.17	0.77
1:A:271:ASN:ND2	1:A:276:PHE:O	2.22	0.72
1:B:131:ASN:ND2	2:D:239:SER:OG	2.23	0.71
1:A:74:ARG:NH2	1:A:297:LYS:O	2.25	0.69
2:C:351:ARG:HB3	6:C:405:MPD:H11	1.74	0.69
1:A:291:CYS:HB2	6:A:406:MPD:H11	1.75	0.68
1:A:20:GLN:NE2	1:A:77:GLU:OE1	2.26	0.68
2:D:219:LYS:NZ	3:D:401:PO4:O1	2.28	0.67
3:B:401:PO4:O1	7:B:501:HOH:O	2.15	0.64
2:C:224:GLN:O	2:C:226:LEU:N	2.28	0.64
2:C:201:ILE:O	2:C:225:ASN:ND2	2.23	0.64
2:D:265:GLU:OE2	2:D:285:ARG:NH2	2.29	0.64
1:B:170:PHE:HE1	1:B:183:MET:HE1	1.63	0.63
2:D:91:ILE:O	2:D:115:ASN:ND2	2.30	0.63
1:B:164:ILE:HD11	1:B:183:MET:HE2	1.82	0.61
2:C:224:GLN:HG3	2:C:225:ASN:H	1.66	0.60
2:D:201:ILE:O	2:D:225:ASN:ND2	2.32	0.60
1:A:246:ARG:HG2	1:A:247:ALA:N	2.16	0.59
2:C:190:MET:HG3	2:C:212:SER:HB2	1.84	0.58
2:C:113:ILE:O	2:C:137:ASN:ND2	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:C:405:MPD:O4	6:C:405:MPD:O2	2.21	0.58
1:B:229:ALA:O	1:B:231:VAL:N	2.37	0.57
2:D:256:GLU:HG2	2:D:278:MET:HB3	1.87	0.56
1:B:74:ARG:NH1	1:B:297:LYS:O	2.39	0.56
2:D:119:LEU:HB3	2:D:122:LEU:HD13	1.87	0.56
1:B:271:ASN:ND2	1:B:277:ASP:HA	2.21	0.55
2:D:260:SER:OG	2:D:280:ASP:OD2	2.24	0.55
2:C:157:ILE:O	2:C:181:ASN:ND2	2.33	0.54
2:D:161:ASP:OD1	2:D:161:ASP:N	2.43	0.52
1:A:113:LYS:HG2	1:A:114:TYR:CE2	2.45	0.52
6:A:406:MPD:O2	6:A:406:MPD:O4	2.26	0.52
2:D:190:MET:HG3	2:D:212:SER:HB2	1.92	0.52
1:B:23:ARG:CZ	1:B:23:ARG:HB3	2.40	0.51
1:B:230:GLU:O	1:B:234:LYS:HD3	2.11	0.51
2:C:143:GLU:HG3	2:C:165:ARG:NH2	2.26	0.51
2:D:124:GLU:HG2	2:D:146:ILE:HD13	1.92	0.51
1:B:174:GLY:O	1:B:246:ARG:NE	2.40	0.50
2:C:289:ILE:HB	2:C:313:ASP:HB3	1.93	0.50
1:A:138:ASP:OD2	1:A:142:ARG:NH2	2.45	0.49
2:C:117:GLU:H	2:C:117:GLU:CD	2.13	0.49
2:C:161:ASP:OD1	2:C:161:ASP:N	2.46	0.49
1:A:131:ASN:ND2	2:C:239:SER:OG	2.46	0.49
1:B:170:PHE:CE1	1:B:183:MET:HE1	2.45	0.49
2:C:256:GLU:HG2	2:C:278:MET:HB3	1.94	0.48
2:D:257:LEU:HD21	2:D:259:LEU:HD11	1.95	0.48
1:A:102:GLU:OE1	1:A:102:GLU:N	2.37	0.48
1:A:138:ASP:OD1	1:A:142:ARG:HD2	2.14	0.48
2:D:123:ARG:HG3	2:D:124:GLU:HG3	1.95	0.48
2:D:221:THR:HG22	2:D:241:ARG:HB2	1.96	0.48
1:B:177:SER:HB2	1:B:203:ASP:HB2	1.96	0.47
2:D:160:VAL:HG21	2:D:182:LEU:HD23	1.96	0.47
2:D:241:ARG:NH2	3:D:401:PO4:O4	2.36	0.47
1:B:216:TRP:CZ3	1:B:227:PHE:HB3	2.49	0.47
1:B:23:ARG:HG2	1:B:24:PRO:HD2	1.96	0.47
2:C:173:ASN:HA	2:C:195:SER:O	2.15	0.47
2:C:94:PHE:O	2:C:119:LEU:HD11	2.14	0.47
1:A:101:LEU:HD12	1:A:143:ARG:HD3	1.97	0.47
1:A:246:ARG:NH2	1:A:249:GLN:OE1	2.48	0.46
1:B:288:THR:O	1:B:288:THR:OG1	2.30	0.46
2:D:321:ARG:HD3	2:D:321:ARG:HA	1.62	0.46
1:B:250:VAL:HG22	1:B:267:PHE:CZ	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:LEU:HD11	1:B:112:ILE:HG22	1.98	0.46
2:D:288:LYS:HG2	2:D:290:GLU:HG3	1.98	0.46
1:B:246:ARG:HD3	1:B:263:LEU:HD21	1.98	0.46
1:A:96:ARG:NH2	1:A:273:CYS:SG	2.89	0.46
2:D:234:VAL:HG22	2:D:256:GLU:HB2	1.97	0.46
2:D:173:ASN:HA	2:D:195:SER:O	2.15	0.45
2:C:120:GLN:NE2	2:C:141:LEU:HA	2.32	0.45
2:C:83:LEU:HB2	2:C:105:LEU:HD23	1.98	0.45
2:C:248:LEU:HD22	2:C:251:LEU:HD12	1.99	0.45
2:D:257:LEU:HD11	2:D:259:LEU:HD21	1.99	0.45
2:D:195:SER:HB2	2:D:217:LYS:HD3	1.99	0.44
2:D:104:CYS:SG	2:D:106:ARG:HG3	2.57	0.44
2:D:310:SER:HG	2:D:312:SER:HG	1.65	0.44
2:D:289:ILE:HB	2:D:313:ASP:HB3	1.98	0.44
1:B:88:LEU:HD11	1:B:121:LEU:HG	2.00	0.44
2:D:77:ASP:CG	2:D:98:LYS:HE2	2.39	0.43
2:C:117:GLU:OE2	2:C:117:GLU:N	2.44	0.43
2:D:224:GLN:HG3	2:D:225:ASN:ND2	2.34	0.43
2:D:237:MET:HB2	2:D:259:LEU:HD23	2.01	0.43
1:A:177:SER:HB2	1:A:203:ASP:HB2	2.01	0.43
1:B:125:HIS:NE2	3:B:401:PO4:O4	2.42	0.42
1:B:267:PHE:CE1	1:B:269:ALA:HB3	2.54	0.42
1:B:194:ASP:OD1	1:B:195:VAL:N	2.43	0.42
2:C:139:GLU:CD	2:C:139:GLU:H	2.22	0.42
2:C:284:ASN:HB3	2:C:285:ARG:H	1.72	0.42
2:C:327:TYR:CD2	2:C:354:ASP:HA	2.55	0.41
1:B:236:LEU:HD21	1:B:244:ILE:HG13	2.01	0.41
2:C:143:GLU:O	2:C:143:GLU:HG2	2.20	0.41
2:D:171:LEU:HB2	2:D:193:LEU:HD23	2.00	0.41
2:D:274:ASN:ND2	2:D:294:HIS:HD2	2.18	0.41
1:B:209:PRO:HA	1:B:227:PHE:O	2.20	0.41
1:A:292:SER:HB3	1:B:294:GLN:HG3	2.02	0.41
2:D:210:LEU:HD23	2:D:210:LEU:HA	1.82	0.41
1:A:194:ASP:OD2	7:A:501:HOH:O	2.22	0.41
1:A:34:GLU:HA	6:A:407:MPD:H51	2.03	0.41
2:C:144:LEU:HD12	2:C:144:LEU:HA	1.88	0.41
1:A:194:ASP:OD1	1:A:195:VAL:N	2.45	0.41
2:D:274:ASN:HD22	2:D:294:HIS:HD2	1.68	0.40
2:D:302:TRP:HA	2:D:327:TYR:HB2	2.03	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	296/305 (97%)	279 (94%)	16 (5%)	1 (0%)	41	66
1	B	289/305 (95%)	268 (93%)	19 (7%)	2 (1%)	22	46
2	C	283/310 (91%)	254 (90%)	28 (10%)	1 (0%)	34	60
2	D	281/310 (91%)	250 (89%)	31 (11%)	0	100	100
All	All	1149/1230 (93%)	1051 (92%)	94 (8%)	4 (0%)	41	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	24	PRO
2	C	225	ASN
1	A	28	VAL
1	B	25	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	A	264/268 (98%)	256 (97%)	8 (3%)	41	70
1	B	254/268 (95%)	247 (97%)	7 (3%)	43	73
2	C	266/291 (91%)	259 (97%)	7 (3%)	46	75
2	D	264/291 (91%)	259 (98%)	5 (2%)	57	82
All	All	1048/1118 (94%)	1021 (97%)	27 (3%)	46	75

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

Mol	Chain	Res	Type
1	A	23	ARG
1	A	96	ARG
1	A	211	LYS
1	A	214	GLN
1	A	234	LYS
1	A	246	ARG
1	A	271	ASN
1	A	297	LYS
1	B	19	VAL
1	B	23	ARG
1	B	26	LYS
1	B	96	ARG
1	B	226	THR
1	B	252	GLU
1	B	288	THR
2	C	84	ASN
2	C	85	HIS
2	C	87	ARG
2	C	117	GLU
2	C	180	GLU
2	C	192	GLU
2	C	205	ASP
2	D	84	ASN
2	D	87	ARG
2	D	125	LEU
2	D	205	ASP
2	D	260	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	ASN
1	A	271	ASN
1	B	27	ASN
1	B	131	ASN
2	C	84	ASN
2	C	115	ASN
2	C	120	GLN
2	C	173	ASN
2	C	184	ASN
2	C	224	GLN

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Mol	Chain	Res	Type
2	C	272	ASN
2	D	274	ASN
2	D	299	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 5 are monoatomic - leaving 12 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	$\# Z > 2$
6	MPD	C	405	-	7,7,7	0.67	0	9,10,10	0.47	0
3	PO4	A	402	-	4,4,4	0.87	0	6,6,6	0.48	0
6	MPD	A	406	-	7,7,7	0.71	0	9,10,10	0.41	0
3	PO4	D	402	-	4,4,4	0.87	0	6,6,6	0.40	0
3	PO4	C	403	-	4,4,4	0.80	0	6,6,6	0.50	0
6	MPD	A	407	-	7,7,7	0.68	0	9,10,10	0.36	0
6	MPD	C	404	-	7,7,7	0.67	0	9,10,10	0.24	0
3	PO4	C	402	-	4,4,4	0.89	0	6,6,6	0.44	0
3	PO4	B	401	4	4,4,4	0.88	0	6,6,6	0.49	0
3	PO4	A	401	4	4,4,4	0.89	0	6,6,6	0.48	0
3	PO4	D	401	-	4,4,4	0.89	0	6,6,6	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PO4	C	401	-	4,4,4	0.90	0	6,6,6	0.44	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
6	MPD	A	407	-	-	0/5/5/5	-
6	MPD	A	406	-	-	0/5/5/5	-
6	MPD	C	404	-	-	0/5/5/5	-
6	MPD	C	405	-	-	4/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	405	MPD	O2-C2-C3-C4
6	C	405	MPD	C1-C2-C3-C4
6	C	405	MPD	CM-C2-C3-C4
6	C	405	MPD	C2-C3-C4-O4

There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	405	MPD	2	0
6	A	406	MPD	2	0
6	A	407	MPD	1	0
3	B	401	PO4	2	0
3	D	401	PO4	2	0

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	298/305 (97%)	-0.18	6 (2%) 65 67	32, 45, 83, 123	0
1	B	291/305 (95%)	-0.06	10 (3%) 45 45	34, 54, 89, 132	0
2	C	285/310 (91%)	-0.33	13 (4%) 32 31	34, 58, 100, 137	0
2	D	283/310 (91%)	0.12	16 (5%) 23 22	39, 73, 117, 138	0
All	All	1157/1230 (94%)	-0.12	45 (3%) 39 38	32, 56, 105, 138	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	97	LEU	6.6
1	A	24	PRO	5.4
2	D	96	VAL	5.2
2	D	100	VAL	5.1
2	D	98	LYS	5.1
2	D	79	GLU	4.5
1	A	275	GLU	3.9
2	C	86	TYR	3.6
2	C	79	GLU	3.6
2	C	78	ALA	3.5
1	B	25	GLY	3.5
2	D	99	LYS	3.3
1	B	237	HIS	3.3
1	B	24	PRO	3.3
2	C	96	VAL	3.2
1	B	23	ARG	3.2
2	C	94	PHE	3.1
1	A	273	CYS	3.1
2	D	321	ARG	3.1
2	C	77	ASP	3.0
1	B	276	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
2	C	85	HIS	2.9
1	A	91	GLY	2.9
2	C	93	GLY	2.9
2	C	360	PHE	2.8
2	D	85	HIS	2.8
2	C	99	LYS	2.7
2	D	77	ASP	2.6
2	D	86	TYR	2.6
1	A	25	GLY	2.6
1	B	277	ASP	2.5
1	B	27	ASN	2.5
2	D	78	ALA	2.5
2	D	164	THR	2.4
2	C	83	LEU	2.4
2	D	186	HIS	2.4
1	A	276	PHE	2.3
2	C	97	LEU	2.2
2	C	87	ARG	2.2
2	D	165	ARG	2.2
2	D	94	PHE	2.2
1	B	245	CYS	2.1
1	B	266	LEU	2.0
2	D	81	VAL	2.0
1	B	274	GLY	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 carbohydrates 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(\AA^2)	Q<0.9
3	PO4	C	403	5/5	0.78	0.52	64,82,106,127	0
6	MPD	A	406	8/8	0.85	0.28	49,68,79,83	0
5	CL	A	405	1/1	0.86	0.24	55,55,55,55	0
6	MPD	C	405	8/8	0.89	0.27	60,74,83,89	0
3	PO4	A	402	5/5	0.90	0.20	73,76,87,107	0
6	MPD	C	404	8/8	0.90	0.40	63,70,83,87	0
3	PO4	D	401	5/5	0.91	0.15	61,82,92,93	0
6	MPD	A	407	8/8	0.92	0.18	49,55,69,69	0
5	CL	A	404	1/1	0.93	0.10	71,71,71,71	0
5	CL	B	403	1/1	0.95	0.06	58,58,58,58	0
3	PO4	D	402	5/5	0.96	0.21	93,94,99,109	0
3	PO4	C	402	5/5	0.97	0.17	71,75,90,94	0
3	PO4	B	401	5/5	0.97	0.20	72,77,82,89	0
4	FE	B	402	1/1	0.97	0.04	66,66,66,66	0
3	PO4	A	401	5/5	0.98	0.09	49,66,72,84	0
3	PO4	C	401	5/5	0.98	0.14	62,67,83,85	0
4	FE	A	403	1/1	0.99	0.05	60,60,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.