



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:44 PM UTC

PDB ID : 1IMD / pdb_00001imd
Title : STRUCTURAL STUDIES OF METAL BINDING BY INOSITOL MONOPHOSPHATASE: EVIDENCE FOR TWO-METAL ION CATALYSIS
Authors : Bone, R.
Deposited on : 1994-02-08
Resolution : 2.60 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

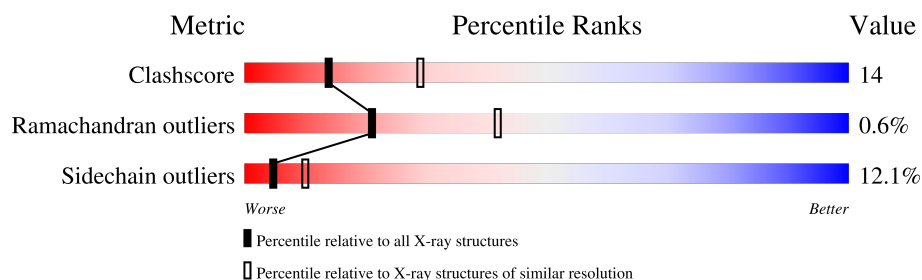
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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))
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	277	
1	B	277	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4157 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.

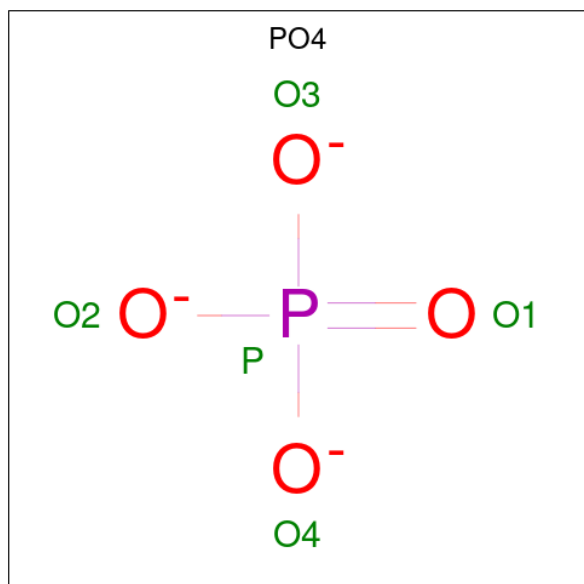
- Molecule 1 is a protein called INOSITOL MONOPHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	53	0	1
			2020	1277	342	383	18			
1	B	273	Total	C	N	O	S	42	0	1
			2073	1307	352	396	18			

- 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	B	2	Total	Mn	0	0
			2	2		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total	O	0	0
			21	21		
4	B	29	Total	O	0	0
			29	29		

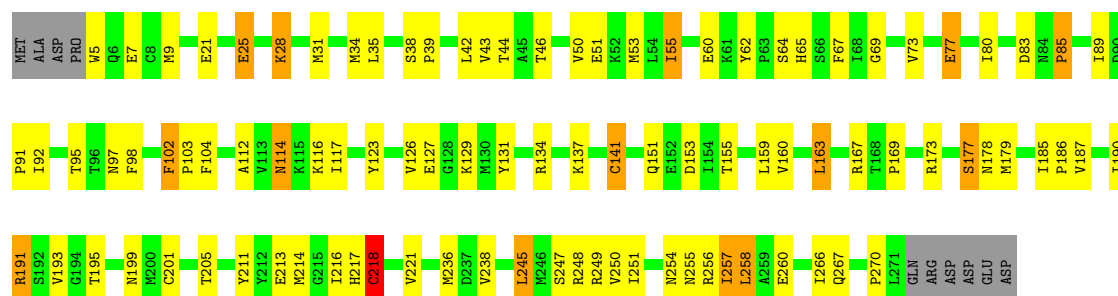
3 Residue-property plots [i](#)

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. 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. 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.

Note EDS was not executed.

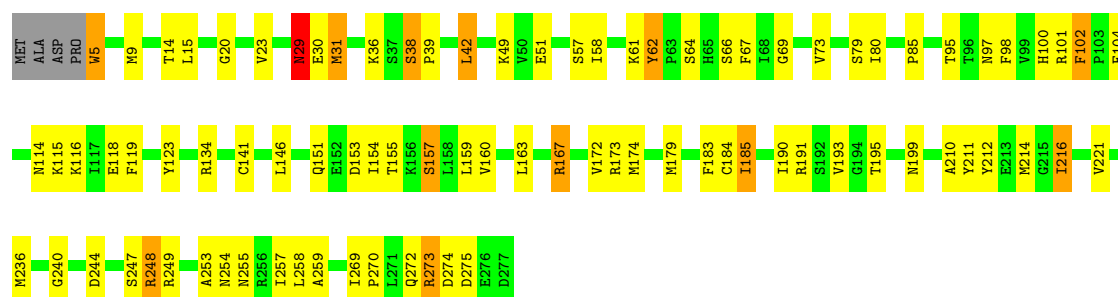
• Molecule 1: INOSITOL MONOPHOSPHATASE

Chain A: 



• Molecule 1: INOSITOL MONOPHOSPHATASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.85Å 85.85Å 153.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.188 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4157	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

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: PO4, MN

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	0.68	1/2053 (0.0%)	0.88	2/2778 (0.1%)
1	B	0.66	0/2106	0.90	4/2849 (0.1%)
All	All	0.67	1/4159 (0.0%)	0.89	6/5627 (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	A	0	3
1	B	0	3
All	All	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	270	PRO	C-N	-5.09	1.26	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	102	PHE	CA-C-N	5.40	126.59	119.84
1	A	102	PHE	C-N-CA	5.40	126.59	119.84
1	B	29	ASN	N-CA-C	5.34	118.52	108.65
1	B	38	SER	CA-C-N	5.07	125.35	119.47
1	B	38	SER	C-N-CA	5.07	125.35	119.47
1	B	185	ILE	CB-CA-C	-5.03	109.02	114.35

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	TYR	Sidechain
1	A	62	TYR	Peptide
1	A	92	ILE	Mainchain
1	B	102	PHE	Peptide
1	B	118	GLU	Mainchain
1	B	212	TYR	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	A	2020	0	2042	63	0
1	B	2073	0	2088	57	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	21	0	0	0	0
4	B	29	0	0	1	0
All	All	4157	0	4130	115	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 14.

All (115) 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:214:MET:HG2	1:B:249:ARG:HG2	1.20	1.13
1:B:236:MET:HE3	1:B:240:GLY:HA2	1.47	0.95
1:A:112:ALA:HB2	1:A:117:ILE:HD13	1.59	0.82
1:A:65:HIS:CD2	1:A:85:PRO:HB2	2.14	0.81
1:A:5:TRP:HB2	1:A:134:ARG:NH2	2.01	0.75
1:A:151:GLN:NE2	1:A:153:ASP:HB3	2.02	0.74
1:A:134:ARG:HB3	1:A:137:LYS:HB2	1.70	0.73
1:B:151:GLN:HE21	1:B:153:ASP:H	1.40	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:ARG:O	1:A:260:GLU:HG3	1.94	0.67
1:B:193:VAL:HG22	1:B:199:ASN:ND2	2.11	0.66
1:B:20:GLY:HA3	1:B:123:TYR:HE2	1.60	0.65
1:A:238:VAL:HG13	1:A:266:ILE:HD11	1.81	0.63
1:B:20:GLY:HA3	1:B:123:TYR:CE2	2.33	0.63
1:B:9:MET:HG3	1:B:119:PHE:CD1	2.33	0.63
1:A:5:TRP:HB2	1:A:134:ARG:HH22	1.63	0.62
1:B:160:VAL:HG23	1:B:211:TYR:HB2	1.81	0.62
1:A:131:TYR:CD1	1:A:141:CYS:HB2	2.34	0.62
1:B:174:MET:HE1	1:B:270:PRO:HD2	1.82	0.62
1:A:214:MET:HG2	1:A:249:ARG:HD2	1.80	0.61
1:B:273:ARG:HD3	1:B:275:ASP:OD1	2.01	0.61
1:B:216:ILE:O	1:B:248:ARG:HD2	2.00	0.61
1:A:151:GLN:HE21	1:A:153:ASP:H	1.48	0.60
1:B:159:LEU:O	1:B:190:ILE:HA	2.02	0.59
1:A:151:GLN:HE21	1:A:153:ASP:HB3	1.68	0.59
1:A:216:ILE:HD13	1:A:221:VAL:CG1	2.33	0.59
1:B:254:ASN:OD1	1:B:258:LEU:HD12	2.02	0.58
1:A:216:ILE:O	1:A:248:ARG:HD2	2.03	0.58
1:A:236:MET:HE1	1:A:256:ARG:HG3	1.86	0.57
1:A:160:VAL:HG23	1:A:211:TYR:HB2	1.86	0.57
1:A:201:CYS:O	1:A:205:THR:HG23	2.04	0.57
1:B:61:LYS:HB3	1:B:62:TYR:CE1	2.39	0.57
1:A:193:VAL:HG22	1:A:199:ASN:OD1	2.04	0.57
1:B:160:VAL:CG2	1:B:211:TYR:HB2	2.35	0.57
1:B:151:GLN:HE21	1:B:153:ASP:N	2.04	0.56
1:B:69:GLY:O	1:B:73:VAL:HG23	2.06	0.55
1:B:154:ILE:O	1:B:157:SER:HB2	2.05	0.55
1:B:51:GLU:OE2	1:B:69:GLY:HA3	2.06	0.55
1:A:255:ASN:OD1	1:A:257:ILE:HB	2.06	0.55
1:A:127:GLU:HB3	1:A:129:LYS:HG3	1.90	0.54
1:B:62:TYR:N	1:B:62:TYR:CD1	2.76	0.53
1:B:248:ARG:HD3	1:B:274:ASP:OD1	2.08	0.53
1:B:193:VAL:HG22	1:B:199:ASN:HD21	1.73	0.53
1:B:100:HIS:O	1:B:101:ARG:HB2	2.07	0.53
1:B:5:TRP:CE3	1:B:5:TRP:HA	2.43	0.53
1:B:38:SER:HB2	1:B:39:PRO:HD2	1.91	0.52
1:A:214:MET:HA	1:A:249:ARG:HG2	1.91	0.52
1:B:9:MET:HG3	1:B:119:PHE:CG	2.45	0.52
1:A:214:MET:HG2	1:A:249:ARG:CD	2.40	0.51
1:A:173:ARG:O	1:A:177:SER:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ARG:NH1	1:B:97:ASN:OD1	2.40	0.51
1:B:159:LEU:CD2	1:B:210:ALA:HB3	2.41	0.51
1:B:5:TRP:HA	1:B:5:TRP:HE3	1.76	0.51
1:A:117:ILE:HD11	1:A:245:LEU:HD21	1.94	0.50
1:A:73:VAL:HA	1:A:77:GLU:O	2.12	0.49
1:A:51:GLU:HB2	1:A:91:PRO:HG3	1.94	0.49
1:A:163:LEU:HB3	1:B:183:PHE:CZ	2.46	0.49
1:A:95:THR:O	1:A:98:PHE:HB3	2.13	0.48
1:A:178:ASN:HD22	1:A:214:MET:HE1	1.78	0.48
1:A:65:HIS:CG	1:A:85:PRO:HB2	2.48	0.48
1:B:36:LYS:HE2	4:B:283:HOH:O	2.14	0.48
1:A:216:ILE:HD13	1:A:221:VAL:HG11	1.95	0.48
1:A:55:ILE:HG13	1:A:67:PHE:CD2	2.48	0.48
1:B:61:LYS:HB3	1:B:62:TYR:CD1	2.49	0.48
1:A:104:PHE:HZ	1:B:104:PHE:HZ	1.61	0.48
1:B:23:VAL:HG13	1:B:98:PHE:CE1	2.49	0.48
1:B:253:ALA:HB3	1:B:259:ALA:HB2	1.95	0.48
1:A:185:ILE:N	1:A:186:PRO:HA	2.29	0.47
1:A:155:THR:HA	1:A:186:PRO:O	2.14	0.47
1:B:29:ASN:HB2	1:B:30:GLU:H	1.52	0.47
1:A:151:GLN:HE21	1:A:153:ASP:N	2.10	0.47
1:B:95:THR:O	1:B:98:PHE:HB3	2.14	0.47
1:A:217:HIS:O	1:A:218:CYS:C	2.58	0.47
1:A:160:VAL:O	1:A:211:TYR:HA	2.14	0.47
1:A:216:ILE:HD13	1:A:221:VAL:HG13	1.96	0.47
1:A:250:VAL:HG22	1:A:251:ILE:N	2.30	0.46
1:A:151:GLN:HE21	1:A:153:ASP:CB	2.28	0.46
1:A:43:VAL:HG22	1:A:44:THR:N	2.31	0.46
1:A:112:ALA:HA	1:A:116:LYS:O	2.17	0.45
1:A:254:ASN:OD1	1:A:258:LEU:HD22	2.15	0.45
1:B:160:VAL:O	1:B:211:TYR:HA	2.16	0.45
1:B:255:ASN:OD1	1:B:257:ILE:HB	2.16	0.45
1:A:103:PRO:HB3	1:A:126:VAL:HG11	1.98	0.45
1:A:247:SER:O	1:A:248:ARG:HB2	2.16	0.45
1:A:178:ASN:ND2	1:A:214:MET:HE1	2.32	0.45
1:B:58:ILE:HB	1:B:67:PHE:CE2	2.52	0.45
1:B:31:MET:HE3	1:B:31:MET:HB2	1.88	0.45
1:B:141:CYS:HB3	1:B:146:LEU:HD21	1.99	0.44
1:B:155:THR:HG22	1:B:185:ILE:HG22	1.99	0.44
1:A:97:ASN:HB3	1:A:102:PHE:O	2.17	0.44
1:A:83:ASP:HB3	1:A:114:ASN:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:ARG:HG2	1:A:260:GLU:HG3	2.00	0.44
1:B:58:ILE:HB	1:B:67:PHE:HE2	1.84	0.43
1:A:69:GLY:O	1:A:73:VAL:HG23	2.19	0.43
1:B:159:LEU:HD22	1:B:210:ALA:HB3	2.00	0.43
1:B:214:MET:HG2	1:B:249:ARG:CG	2.14	0.43
1:A:5:TRP:HA	1:A:5:TRP:CE3	2.54	0.42
1:B:269:ILE:HA	1:B:270:PRO:HD2	1.80	0.42
1:B:42:LEU:HD23	1:B:42:LEU:N	2.34	0.42
1:A:236:MET:HB2	1:A:236:MET:HE3	1.72	0.42
1:A:159:LEU:O	1:A:190:ILE:HA	2.20	0.42
1:A:191:ARG:HG2	1:B:102:PHE:CZ	2.54	0.42
1:A:55:ILE:HD12	1:A:89:ILE:HD13	2.01	0.42
1:B:5:TRP:HB2	1:B:134:ARG:NH2	2.34	0.42
1:A:187:VAL:O	1:B:167:ARG:NH1	2.52	0.42
1:B:258:LEU:HD23	1:B:258:LEU:HA	1.91	0.42
1:B:216:ILE:HD13	1:B:221:VAL:HG11	2.01	0.41
1:A:131:TYR:CE1	1:A:141:CYS:HB2	2.55	0.41
1:B:244:ASP:HB3	1:B:247:SER:HB3	2.02	0.41
1:B:114:ASN:O	1:B:115:LYS:HB2	2.20	0.41
1:B:163:LEU:CD2	1:B:179:MET:HE1	2.50	0.41
1:A:25:GLU:O	1:A:28:LYS:HG2	2.21	0.41
1:A:35:LEU:HD13	1:A:39:PRO:HA	2.01	0.41
1:B:216:ILE:HD13	1:B:221:VAL:CG1	2.51	0.41
1:A:38:SER:HB2	1:A:39:PRO:HD2	2.03	0.41
1:A:46:THR:O	1:A:50:VAL:HG23	2.21	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	265/277 (96%)	248 (94%)	15 (6%)	2 (1%)	16	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	271/277 (98%)	254 (94%)	16 (6%)	1 (0%)	30	51
All	All	536/554 (97%)	502 (94%)	31 (6%)	3 (1%)	21	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	218	CYS
1	B	85	PRO
1	A	85	PRO

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	221/231 (96%)	192 (87%)	29 (13%)	4	8
1	B	227/231 (98%)	202 (89%)	25 (11%)	6	13
All	All	448/462 (97%)	394 (88%)	54 (12%)	5	10

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	9	MET
1	A	21	GLU
1	A	25	GLU
1	A	28	LYS
1	A	31	MET
1	A	34	MET
1	A	42	LEU
1	A	53	MET
1	A	55	ILE
1	A	60	GLU
1	A	64	SER
1	A	77	GLU

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Mol	Chain	Res	Type
1	A	80	ILE
1	A	114	ASN
1	A	141	CYS
1	A	163	LEU
1	A	167	ARG
1	A	169	PRO
1	A	177	SER
1	A	179	MET
1	A	191	ARG
1	A	195	THR
1	A	213	GLU
1	A	218	CYS
1	A	245	LEU
1	A	257	ILE
1	A	258	LEU
1	A	267	GLN
1	B	5	TRP
1	B	14	THR
1	B	15	LEU
1	B	29	ASN
1	B	31	MET
1	B	42	LEU
1	B	49	LYS
1	B	57	SER
1	B	62	TYR
1	B	64	SER
1	B	66	SER
1	B	79	SER
1	B	80	ILE
1	B	116	LYS
1	B	157	SER
1	B	167	ARG
1	B	172	VAL
1	B	173	ARG
1	B	184	CYS
1	B	191	ARG
1	B	195	THR
1	B	216	ILE
1	B	248	ARG
1	B	272	GLN
1	B	273	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	48	GLN
1	A	147	GLN
1	A	151	GLN
1	B	32	ASN
1	B	48	GLN
1	B	151	GLN
1	B	199	ASN
1	B	267	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 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	PO4	A	280	2	4,4,4	0.79	0	6,6,6	0.37	0
3	PO4	B	280	2	4,4,4	0.64	0	6,6,6	0.58	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 [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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.