



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:01 PM GMT

PDB ID : 1DPM  
Title : THREE-DIMENSIONAL STRUCTURE OF THE ZINC-CONTAINING PHOSPHOTRIESTERASE WITH BOUND SUBSTRATE ANALOG DI-ETHYL 4-METHYLBENZYLPHOSPHONATE  
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Deposited on : 1996-02-13  
Resolution : 2.10 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

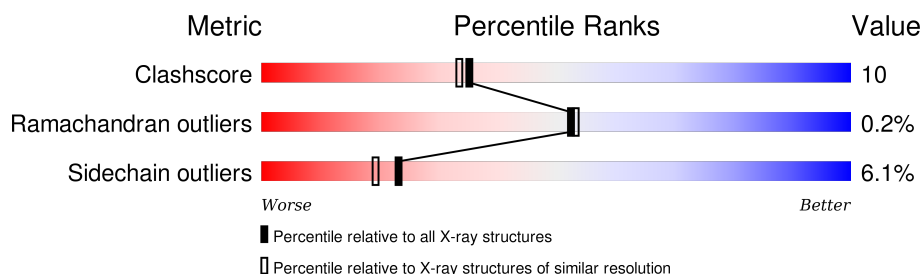
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	329	
1	B	329	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5361 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.

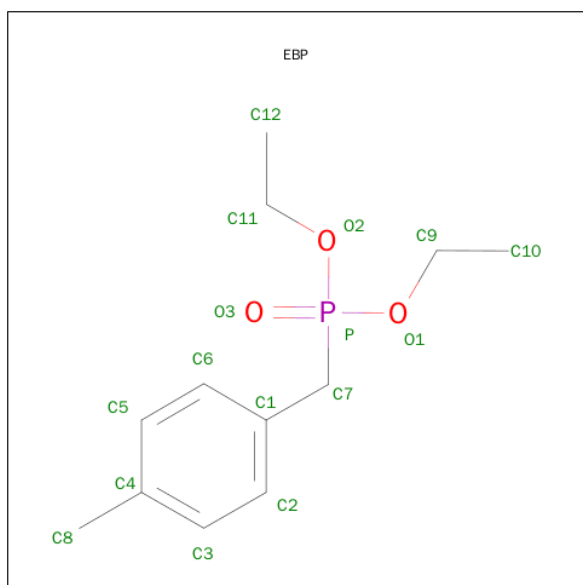
- Molecule 1 is a protein called PHOSPHOTRIESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2515	1592	449	467	7			
1	B	329	Total	C	N	O	S	0	0	0
			2509	1589	446	467	7			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

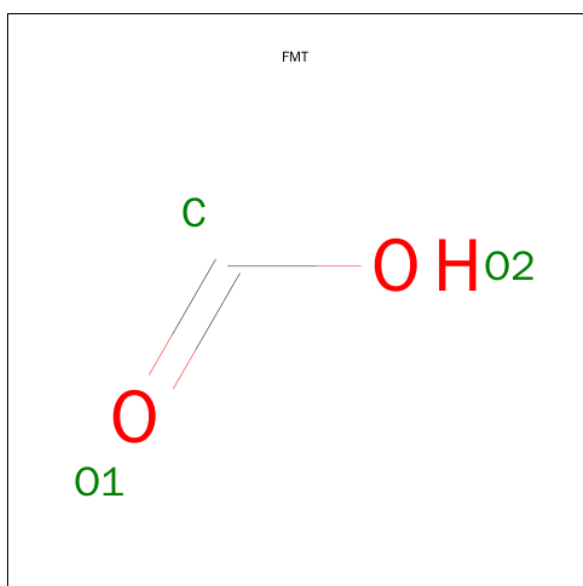
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is DIETHYL 4-METHYLBENZYLPHOSPHONATE (three-letter code: EBP) (formula: C<sub>12</sub>H<sub>19</sub>O<sub>3</sub>P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			16	12	3	1		
3	A	1	Total	C	O	P	0	0
			16	12	3	1		
3	B	1	Total	C	O	P	0	0
			16	12	3	1		
3	B	1	Total	C	O	P	0	0
			16	12	3	1		

- Molecule 4 is FORMIC ACID (three-letter code: FMT) (formula: CH<sub>2</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	120	Total	O	0	0
			120	120		
5	B	143	Total	O	0	0
			143	143		

### 3 Residue-property plots [i](#)

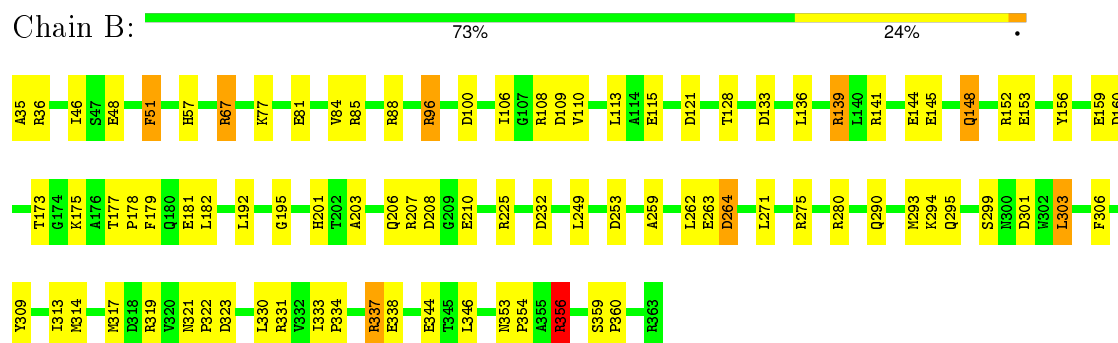
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

#### • Molecule 1: PHOSPHOTRIESTERASE



#### • Molecule 1: PHOSPHOTRIESTERASE



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	129.60 Å 91.40 Å 69.40 Å 90.00° 91.90° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.10)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	TNT	Depositor
R, $R_{free}$	0.154 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5361	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, EBP, FMT

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	1.07	10/2563 (0.4%)	1.37	33/3479 (0.9%)
1	B	1.05	11/2557 (0.4%)	1.37	35/3472 (1.0%)
All	All	1.06	21/5120 (0.4%)	1.37	68/6951 (1.0%)

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	B	1	0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	118	ARG	CD-NE	9.82	1.63	1.46
1	A	263	GLU	CD-OE1	7.95	1.34	1.25
1	B	344	GLU	CD-OE2	7.52	1.33	1.25
1	A	338	GLU	CD-OE2	7.23	1.33	1.25
1	B	144	GLU	CD-OE1	6.95	1.33	1.25
1	B	115	GLU	CD-OE1	6.80	1.33	1.25
1	B	48	GLU	CD-OE2	6.68	1.32	1.25
1	A	344	GLU	CD-OE1	6.66	1.32	1.25
1	B	210	GLU	CD-OE1	6.64	1.32	1.25
1	B	263	GLU	CD-OE1	6.57	1.32	1.25
1	A	115	GLU	CD-OE1	6.49	1.32	1.25
1	B	181	GLU	CD-OE1	6.44	1.32	1.25
1	A	210	GLU	CD-OE1	6.27	1.32	1.25
1	B	338	GLU	CD-OE2	5.92	1.32	1.25
1	B	81	GLU	CD-OE1	5.89	1.32	1.25
1	A	145	GLU	CD-OE1	5.88	1.32	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	81	GLU	CD-OE1	5.80	1.32	1.25
1	A	181	GLU	CD-OE1	5.66	1.31	1.25
1	A	48	GLU	CD-OE2	5.60	1.31	1.25
1	B	159	GLU	CD-OE1	5.52	1.31	1.25
1	B	153	GLU	CD-OE1	5.24	1.31	1.25

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	139	ARG	NE-CZ-NH1	13.04	126.82	120.30
1	A	139	ARG	NE-CZ-NH1	10.21	125.40	120.30
1	A	356	ARG	NE-CZ-NH1	9.46	125.03	120.30
1	A	85	ARG	NE-CZ-NH2	-8.88	115.86	120.30
1	B	141	ARG	NE-CZ-NH1	8.78	124.69	120.30
1	A	232	ASP	CB-CG-OD1	-8.18	110.94	118.30
1	A	164	ARG	NE-CZ-NH1	8.07	124.34	120.30
1	A	208	ASP	CB-CG-OD1	-7.87	111.22	118.30
1	A	139	ARG	NE-CZ-NH2	-7.66	116.47	120.30
1	A	133	ASP	CB-CG-OD1	-7.61	111.45	118.30
1	B	139	ARG	NE-CZ-NH2	-7.53	116.54	120.30
1	B	323	ASP	CB-CG-OD1	-7.39	111.65	118.30
1	A	253	ASP	CB-CG-OD2	7.39	124.95	118.30
1	B	323	ASP	CB-CG-OD2	7.37	124.94	118.30
1	B	301	ASP	CB-CG-OD2	7.33	124.90	118.30
1	A	232	ASP	CB-CG-OD2	6.98	124.58	118.30
1	A	235	ASP	CB-CG-OD1	-6.86	112.13	118.30
1	B	108	ARG	NE-CZ-NH1	6.84	123.72	120.30
1	B	133	ASP	CB-CG-OD1	-6.83	112.15	118.30
1	B	232	ASP	CB-CG-OD2	6.83	124.45	118.30
1	B	253	ASP	CB-CG-OD2	6.82	124.44	118.30
1	B	253	ASP	CB-CG-OD1	-6.81	112.17	118.30
1	B	109	ASP	CB-CG-OD1	6.76	124.39	118.30
1	B	121	ASP	CB-CG-OD2	-6.67	112.30	118.30
1	B	280	ARG	NE-CZ-NH1	6.57	123.58	120.30
1	A	280	ARG	NE-CZ-NH1	6.55	123.57	120.30
1	B	152	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	A	133	ASP	CB-CG-OD2	6.51	124.16	118.30
1	B	232	ASP	CB-CG-OD1	-6.49	112.46	118.30
1	A	105	ASP	CB-CG-OD1	-6.45	112.50	118.30
1	B	121	ASP	CB-CG-OD1	6.43	124.09	118.30
1	A	109	ASP	CB-CG-OD1	6.38	124.04	118.30
1	B	100	ASP	CB-CG-OD2	6.36	124.02	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	85	ARG	NE-CZ-NH2	-6.32	117.14	120.30
1	A	264	ASP	CB-CG-OD1	-6.26	112.67	118.30
1	A	315	ASP	CB-CG-OD2	-6.26	112.67	118.30
1	A	301	ASP	CB-CG-OD2	6.17	123.86	118.30
1	B	263	GLU	N-CA-CB	6.17	121.71	110.60
1	A	208	ASP	CB-CG-OD2	6.13	123.82	118.30
1	A	356	ARG	NE-CZ-NH2	-6.12	117.24	120.30
1	A	85	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	B	100	ASP	CB-CG-OD1	-5.93	112.96	118.30
1	B	264	ASP	CB-CG-OD1	5.92	123.63	118.30
1	B	160	ASP	CB-CG-OD2	5.92	123.62	118.30
1	B	85	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	A	203	ALA	N-CA-C	-5.83	95.25	111.00
1	A	235	ASP	CB-CG-OD2	5.75	123.48	118.30
1	B	225	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	A	152	ARG	NE-CZ-NH2	-5.73	117.43	120.30
1	A	100	ASP	CB-CG-OD1	-5.72	113.15	118.30
1	A	264	ASP	CB-CG-OD2	5.71	123.44	118.30
1	B	108	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	B	264	ASP	CB-CG-OD2	-5.67	113.20	118.30
1	B	133	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	76	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	A	225	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	B	160	ASP	CB-CG-OD1	-5.52	113.33	118.30
1	A	108	ARG	NE-CZ-NH2	-5.37	117.61	120.30
1	A	100	ASP	CB-CG-OD2	5.35	123.12	118.30
1	A	239	TYR	CB-CG-CD2	-5.35	117.79	121.00
1	A	121	ASP	CB-CG-OD1	5.28	123.05	118.30
1	B	203	ALA	N-CA-C	-5.23	96.87	111.00
1	B	356	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	B	208	ASP	CB-CG-OD1	-5.22	113.61	118.30
1	B	337	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	B	306	PHE	N-CA-CB	-5.16	101.31	110.60
1	A	236	ASP	CB-CG-OD1	-5.08	113.73	118.30
1	B	128	THR	N-CA-CB	5.01	119.81	110.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	263	GLU	CA

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	2515	0	2542	56	0
1	B	2509	0	2531	43	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	32	0	38	7	0
3	B	32	0	38	10	0
4	A	3	0	0	1	0
4	B	3	0	0	1	0
5	A	120	0	0	3	0
5	B	143	0	0	6	0
All	All	5361	0	5149	100	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 10.

All (100) 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:156:TYR:CG	3:B:903:EBP:H122	2.16	0.80
1:B:271:LEU:HD21	3:B:902:EBP:H5	1.70	0.73
1:A:333:ILE:HB	1:A:334:PRO:HD3	1.72	0.70
1:B:156:TYR:CD2	3:B:903:EBP:H122	2.27	0.70
1:B:293:MET:HG2	5:B:997:HOH:O	1.93	0.67
1:B:353:ASN:HB2	1:B:354:PRO:HD3	1.78	0.66
1:B:96:ARG:HD3	5:B:1032:HOH:O	1.96	0.63
1:A:46:ILE:HD12	1:A:359:SER:OG	1.98	0.63
1:A:271:LEU:CD2	3:A:900:EBP:H6	2.28	0.63
1:B:317:MET:SD	3:B:902:EBP:H82	2.39	0.62
1:A:303:LEU:HD11	1:A:317:MET:HE1	1.81	0.62
1:A:271:LEU:HD21	3:A:900:EBP:H6	1.82	0.61
1:B:156:TYR:CD1	3:B:903:EBP:H122	2.36	0.61
1:B:145:GLU:O	1:B:148:GLN:HB2	2.01	0.61
1:A:303:LEU:CD1	1:A:317:MET:HE1	2.31	0.60
1:A:312:ASN:O	1:A:315:ASP:HB2	2.01	0.60
1:B:57:HIS:O	1:B:303:LEU:HA	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:VAL:HG22	1:A:163:ILE:HD13	1.84	0.59
1:B:294:LYS:O	1:B:356:ARG:NH2	2.36	0.57
1:B:173:THR:HG23	1:B:173:THR:O	2.04	0.56
1:B:271:LEU:HD21	3:B:902:EBP:C5	2.35	0.56
1:B:333:ILE:HG23	1:B:346:LEU:HD13	1.87	0.55
3:B:902:EBP:H72	5:B:924:HOH:O	2.07	0.54
1:A:314:MET:CE	1:A:317:MET:HE2	2.38	0.54
1:B:317:MET:SD	3:B:902:EBP:C8	2.96	0.53
1:A:57:HIS:O	1:A:303:LEU:HA	2.07	0.53
1:A:353:ASN:HB2	1:A:354:PRO:HD3	1.89	0.53
1:B:271:LEU:CD2	3:B:902:EBP:H5	2.40	0.52
1:B:206:GLN:O	1:B:207:ARG:HB2	2.09	0.52
1:B:275:ARG:HG3	5:B:1000:HOH:O	2.11	0.51
1:B:46:ILE:HD12	1:B:359:SER:OG	2.11	0.50
1:A:314:MET:HE3	1:A:317:MET:CE	2.40	0.50
1:A:236:ASP:OD1	1:A:238:SER:OG	2.26	0.50
1:B:84:VAL:O	1:B:88:ARG:HG3	2.12	0.50
1:A:55:HIS:CE1	1:A:230:HIS:CE1	3.00	0.50
1:A:255:ILE:HB	1:A:256:PRO:HD3	1.94	0.49
1:A:114:ALA:O	1:A:118:ARG:HG3	2.12	0.49
1:A:230:HIS:O	1:A:233:ASP:HB2	2.12	0.49
1:A:201:HIS:HB2	4:A:902:FMT:O1	2.13	0.49
1:B:271:LEU:CD2	1:B:317:MET:HE1	2.43	0.48
1:A:50:GLY:HA3	1:A:96:ARG:HD3	1.94	0.48
1:A:314:MET:CE	1:A:317:MET:CE	2.92	0.48
1:A:271:LEU:HD21	3:A:900:EBP:C6	2.43	0.48
1:A:317:MET:SD	3:A:900:EBP:C8	3.02	0.48
1:B:136:LEU:CD2	1:B:139:ARG:NH2	2.77	0.48
1:B:271:LEU:HD21	1:B:317:MET:HE1	1.96	0.47
1:B:35:ALA:N	5:B:936:HOH:O	2.47	0.47
1:B:353:ASN:N	1:B:354:PRO:CD	2.78	0.47
1:A:317:MET:HB3	1:A:317:MET:HE2	1.60	0.47
1:B:309:TYR:HD2	1:B:313:ILE:HG12	1.80	0.47
1:B:353:ASN:N	1:B:354:PRO:HD2	2.30	0.47
1:A:317:MET:SD	3:A:900:EBP:H82	2.55	0.46
1:B:201:HIS:HB2	4:B:904:FMT:O2	2.15	0.46
1:A:67:ARG:CZ	1:B:67:ARG:HH12	2.28	0.46
1:B:321:ASN:HA	1:B:322:PRO:HD2	1.61	0.46
1:B:259:ALA:O	1:B:262:LEU:HB2	2.16	0.46
1:A:333:ILE:HG23	1:A:346:LEU:HD13	1.98	0.46
1:B:179:PHE:O	1:B:182:LEU:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:ILE:HG22	1:B:106:ILE:O	2.16	0.46
1:A:335:PHE:O	1:A:339:LYS:HG2	2.17	0.45
1:A:117:SER:HB2	1:A:124:ILE:HD12	1.99	0.44
1:A:110:VAL:HG22	1:A:163:ILE:CD1	2.46	0.44
1:A:133:ASP:N	1:A:134:PRO:CD	2.81	0.44
1:A:202:THR:HB	1:A:208:ASP:HB2	2.00	0.44
1:A:51:PHE:C	1:A:51:PHE:CD1	2.92	0.44
1:A:67:ARG:NH1	1:B:67:ARG:NH1	2.66	0.44
3:A:901:EBP:H6	3:A:901:EBP:H92	2.00	0.44
1:A:51:PHE:C	1:A:51:PHE:HD1	2.22	0.43
1:A:322:PRO:HD2	5:A:1003:HOH:O	2.18	0.43
1:A:301:ASP:O	3:A:900:EBP:H3	2.19	0.43
1:B:195:GLY:O	1:B:360:PRO:HA	2.18	0.43
1:A:329:PRO:HA	1:A:333:ILE:HG12	2.00	0.43
1:A:295:GLN:HB3	1:A:295:GLN:HE21	1.68	0.43
1:B:177:THR:HB	1:B:178:PRO:HD2	2.01	0.42
1:B:271:LEU:CD2	1:B:317:MET:CE	2.97	0.42
1:B:51:PHE:C	1:B:51:PHE:CD1	2.92	0.42
1:B:51:PHE:C	1:B:51:PHE:HD1	2.21	0.42
1:A:173:THR:HG23	1:A:173:THR:O	2.20	0.42
1:B:192:LEU:HA	1:B:192:LEU:HD23	1.83	0.42
1:A:358:LEU:HD23	1:A:358:LEU:HA	1.89	0.42
1:A:176:ALA:N	1:A:211:GLN:OE1	2.47	0.42
1:A:259:ALA:O	1:A:262:LEU:HB2	2.20	0.42
1:A:55:HIS:CE1	1:A:230:HIS:HE1	2.38	0.42
1:A:206:GLN:O	1:A:207:ARG:HB2	2.20	0.42
3:B:903:EBP:H71	3:B:903:EBP:H91	1.87	0.41
1:B:330:LEU:O	1:B:334:PRO:HG2	2.20	0.41
1:A:326:ALA:O	1:A:329:PRO:HD2	2.20	0.41
1:A:239:TYR:O	1:A:242:ALA:HB3	2.20	0.41
1:A:82:LYS:NZ	5:A:921:HOH:O	2.53	0.41
1:B:333:ILE:HG22	5:B:931:HOH:O	2.19	0.41
1:B:57:HIS:HB2	1:B:303:LEU:HB3	2.02	0.41
1:A:237:LEU:HD12	1:A:240:LEU:HD12	2.03	0.41
1:A:339:LYS:HA	1:A:339:LYS:HD3	1.89	0.41
1:A:143:VAL:HG12	5:A:928:HOH:O	2.21	0.41
1:A:321:ASN:HA	1:A:322:PRO:HD2	1.80	0.41
1:A:195:GLY:HA3	1:A:362:LEU:HD13	2.03	0.41
1:A:105:ASP:OD2	1:A:131:TRP:HB3	2.21	0.41
1:A:73:PHE:CZ	1:A:307:SER:HB3	2.56	0.40
1:A:71:GLU:CD	1:A:71:GLU:H	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:ALA:HB2	1:A:272:LEU:HB2	2.03	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	327/329 (99%)	313 (96%)	13 (4%)	1 (0%)	46	45
1	B	327/329 (99%)	318 (97%)	9 (3%)	0	100	100
All	All	654/658 (99%)	631 (96%)	22 (3%)	1 (0%)	52	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	101	VAL

### 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	264/264 (100%)	252 (96%)	12 (4%)	34	32
1	B	263/264 (100%)	243 (92%)	20 (8%)	16	12
All	All	527/528 (100%)	495 (94%)	32 (6%)	23	19

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	PHE
1	A	67	ARG
1	A	211	GLN
1	A	264	ASP
1	A	271	LEU
1	A	299	SER
1	A	303	LEU
1	A	306	PHE
1	A	314	MET
1	A	337	ARG
1	A	344	GLU
1	A	362	LEU
1	B	36	ARG
1	B	51	PHE
1	B	67	ARG
1	B	77	LYS
1	B	96	ARG
1	B	110	VAL
1	B	113	LEU
1	B	148	GLN
1	B	175	LYS
1	B	249	LEU
1	B	264	ASP
1	B	290	GLN
1	B	295	GLN
1	B	299	SER
1	B	303	LEU
1	B	314	MET
1	B	319	ARG
1	B	331	ARG
1	B	337	ARG
1	B	356	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	212	GLN
1	A	295	GLN
1	B	212	GLN
1	B	290	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	EBP	A	900	-	16,16,16	3.93	3 (18%)	21,21,21	1.34	1 (4%)
3	EBP	A	901	-	16,16,16	4.02	3 (18%)	21,21,21	0.84	1 (4%)
4	FMT	A	902	1,2	0,2,2	0.00	-	0,1,1	0.00	-
3	EBP	B	902	-	16,16,16	3.85	3 (18%)	21,21,21	1.71	1 (4%)
3	EBP	B	903	-	16,16,16	4.19	4 (25%)	21,21,21	1.32	1 (4%)
4	FMT	B	904	1,2	0,2,2	0.00	-	0,1,1	0.00	-

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	EBP	A	900	-	-	0/13/13/13	0/1/1/1
3	EBP	A	901	-	-	0/13/13/13	0/1/1/1
4	FMT	A	902	1,2	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EBP	B	902	-	-	0/13/13/13	0/1/1/1
3	EBP	B	903	-	-	0/13/13/13	0/1/1/1
4	FMT	B	904	1,2	-	0/0/0/0	0/0/0/0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	903	EBP	P-C7	-15.81	1.60	1.79
3	A	901	EBP	P-C7	-15.21	1.61	1.79
3	A	900	EBP	P-C7	-14.82	1.61	1.79
3	B	902	EBP	P-C7	-14.17	1.62	1.79
3	B	903	EBP	C2-C1	2.36	1.43	1.38
3	B	903	EBP	C6-C5	2.54	1.43	1.38
3	A	900	EBP	C6-C5	2.72	1.43	1.38
3	A	900	EBP	C3-C2	2.78	1.43	1.38
3	A	901	EBP	C6-C5	2.97	1.44	1.38
3	B	903	EBP	C3-C2	3.29	1.44	1.38
3	A	901	EBP	C3-C2	3.45	1.45	1.38
3	B	902	EBP	C3-C2	3.45	1.45	1.38
3	B	902	EBP	C6-C5	3.63	1.45	1.38

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	EBP	C2-C3-C4	-2.65	117.82	121.39
3	B	903	EBP	P-C7-C1	4.77	124.58	113.84
3	A	900	EBP	P-C7-C1	4.79	124.62	113.84
3	B	902	EBP	P-C7-C1	7.35	130.40	113.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	900	EBP	6	0
3	A	901	EBP	1	0
4	A	902	FMT	1	0
3	B	902	EBP	6	0
3	B	903	EBP	4	0
4	B	904	FMT	1	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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.