



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:38 PM GMT

PDB ID : 1L8P  
Title : Mg-phosphonoacetohydroxamate complex of S39A yeast enolase 1  
Authors : Poyner, R.R.; Larsen, T.M.; Wong, S.W.; Reed, G.H.  
Deposited on : 2002-03-21  
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

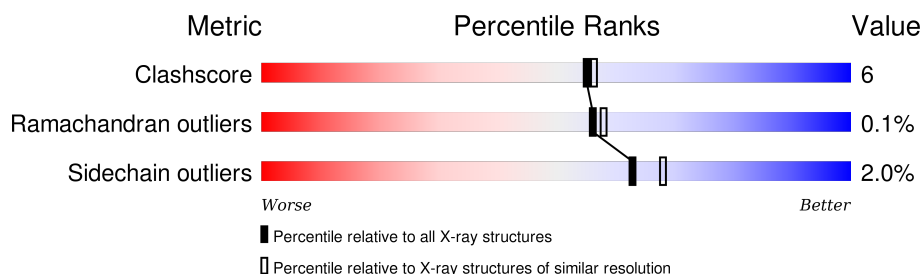
# 1 Overall quality at a glance

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	436	 81% 19%
1	B	436	 88% 12%
1	C	436	 83% 16% .
1	D	436	 86% 14%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 14636 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 enolase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	436	Total	C	N	O	S	0	0	0
			3290	2079	570	635	6			
1	B	436	Total	C	N	O	S	0	0	0
			3290	2079	570	635	6			
1	C	436	Total	C	N	O	S	0	0	0
			3290	2079	570	635	6			
1	D	436	Total	C	N	O	S	0	0	0
			3290	2079	570	635	6			

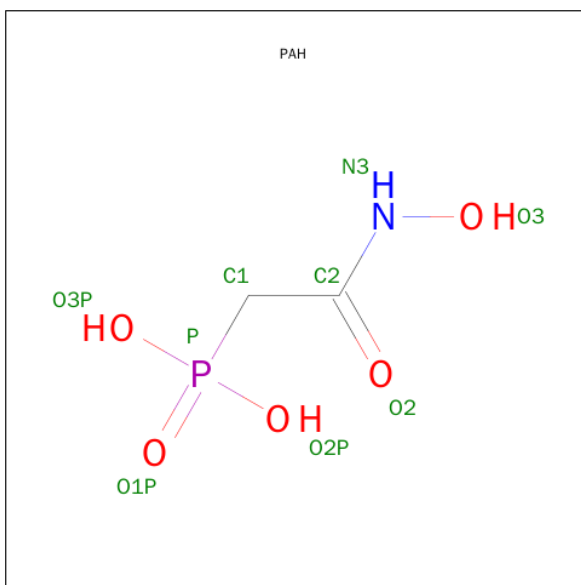
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	ALA	SER	ENGINEERED	UNP P00924
B	539	ALA	SER	ENGINEERED	UNP P00924
C	1039	ALA	SER	ENGINEERED	UNP P00924
D	1539	ALA	SER	ENGINEERED	UNP P00924

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		
2	D	2	Total	Mg	0	0
			2	2		
2	C	2	Total	Mg	0	0
			2	2		

- Molecule 3 is PHOSPHONOACETOHYDROXAMIC ACID (three-letter code: PAH) (formula: C<sub>2</sub>H<sub>6</sub>NO<sub>5</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			9	2	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			9	2	1	5	1		
3	C	1	Total	C	N	O	P	0	0
			9	2	1	5	1		
3	D	1	Total	C	N	O	P	0	0
			9	2	1	5	1		

- Molecule 4 is water.

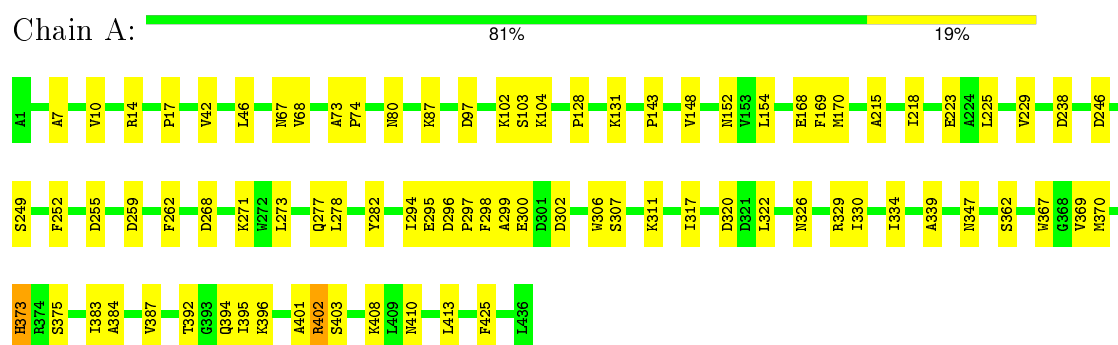
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	339	Total	O	0	0
			339	339		
4	B	388	Total	O	0	0
			388	388		
4	C	328	Total	O	0	0
			328	328		
4	D	377	Total	O	0	0
			377	377		

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

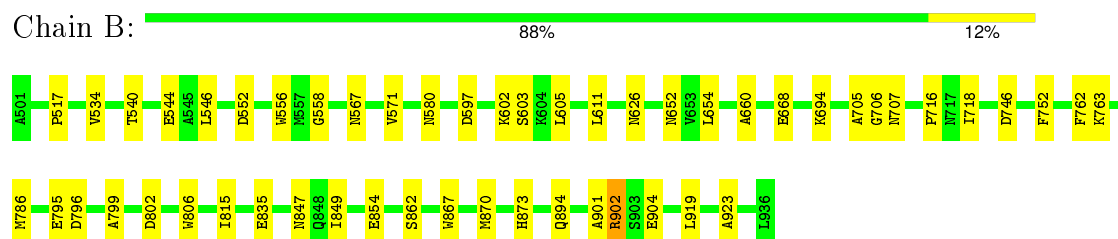
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

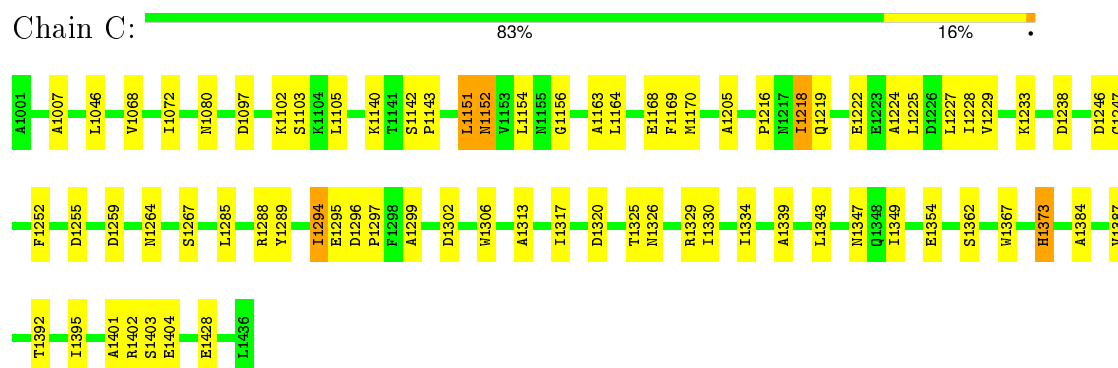
- Molecule 1: enolase 1



- Molecule 1: enolase 1



- Molecule 1: enolase 1

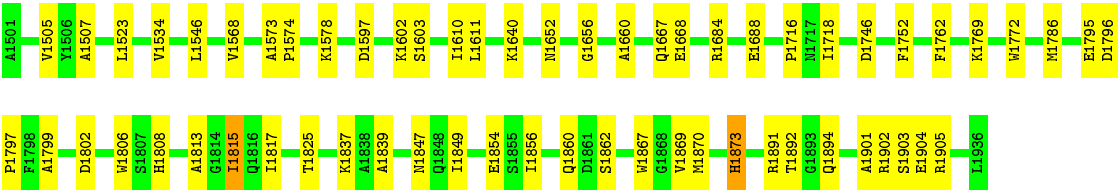


- Molecule 1: enolase 1

Chain D: 

86%

14%



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	63.20 Å   164.10 Å   84.60 Å 90.00°   90.40°   90.00°	Depositor
Resolution (Å)	20.00 – 2.10	Depositor
% Data completeness (in resolution range)	96.7 (20.00-2.10)	Depositor
$R_{merge}$	0.03	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.181 , 0.223	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	14636	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	15.0	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: PAH, MG

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	0/3350	0.57	0/4533
1	B	0.32	0/3350	0.58	0/4533
1	C	0.32	0/3350	0.56	0/4533
1	D	0.33	0/3350	0.58	0/4533
All	All	0.32	0/13400	0.57	0/18132

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 [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	3290	0	3301	52	0
1	B	3290	0	3298	30	0
1	C	3290	0	3298	51	0
1	D	3290	0	3298	40	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	9	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	9	0	3	0	0
3	C	9	0	3	0	0
3	D	9	0	3	0	0
4	A	339	0	0	4	0
4	B	388	0	0	2	0
4	C	328	0	0	1	0
4	D	377	0	0	3	0
All	All	14636	0	13207	165	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 6.

All (165) 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:C:1168:GLU:HB2	1:C:1246:ASP:HB3	1.59	0.85
1:A:168:GLU:HB2	1:A:246:ASP:HB3	1.66	0.76
1:A:330:ILE:O	1:A:334:ILE:HG12	1.89	0.73
1:C:1362:SER:O	1:C:1367:TRP:HB2	1.91	0.70
1:C:1072:ILE:HD11	1:C:1105:LEU:HD13	1.75	0.69
1:B:668:GLU:HB2	1:B:746:ASP:HB3	1.77	0.67
1:C:1218:ILE:HD12	1:C:1224:ALA:HB2	1.78	0.65
1:B:799:ALA:HB3	1:B:802:ASP:HB2	1.80	0.63
1:A:7:ALA:HB2	1:A:68:VAL:HG11	1.81	0.62
1:C:1163:ALA:HB3	1:C:1219:GLN:HA	1.82	0.62
1:C:1246:ASP:HA	1:C:1295:GLU:HB3	1.81	0.61
1:C:1205:ALA:HB1	1:C:1216:PRO:HB3	1.82	0.61
1:D:1799:ALA:HB3	1:D:1802:ASP:HB2	1.83	0.60
1:C:1330:ILE:O	1:C:1334:ILE:HG12	2.02	0.60
1:D:1815:ILE:H	1:D:1815:ILE:HD13	1.66	0.60
1:C:1299:ALA:HB3	1:C:1302:ASP:HB2	1.84	0.59
1:D:1668:GLU:HB2	1:D:1746:ASP:HB3	1.84	0.59
1:D:1746:ASP:HA	1:D:1795:GLU:HB3	1.84	0.59
1:A:296:ASP:OD2	1:A:320:ASP:HB3	2.05	0.57
1:A:298:PHE:HB2	1:A:306:TRP:CD1	2.41	0.56
1:A:154:LEU:HB3	1:A:169:PHE:HB2	1.87	0.55
1:C:1297:PRO:HD2	1:C:1306:TRP:CH2	2.40	0.55
1:A:97:ASP:O	1:A:104:LYS:HD2	2.06	0.55
1:C:1401:ALA:O	1:C:1402:ARG:HB2	2.06	0.55
1:C:1387:VAL:HG21	1:C:1395:ILE:HB	1.87	0.55
1:D:1505:VAL:HG22	1:D:1523:LEU:HD12	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1152:ASN:OD1	1:C:1168:GLU:HG2	2.07	0.55
1:A:299:ALA:HB3	1:A:302:ASP:HB2	1.90	0.54
1:C:1140:LYS:HG2	1:C:1142:SER:H	1.72	0.54
1:C:1384:ALA:O	1:C:1387:VAL:HG12	2.08	0.54
1:A:104:LYS:HE3	4:A:2589:HOH:O	2.07	0.54
1:D:1849:ILE:HG12	1:D:1854:GLU:HB3	1.89	0.54
1:B:716:PRO:HG2	1:B:718:ILE:HG13	1.89	0.53
1:B:746:ASP:HA	1:B:795:GLU:HB3	1.88	0.53
1:A:170:MET:SD	1:A:396:LYS:HD2	2.49	0.53
1:B:567:ASN:O	1:B:571:VAL:HB	2.09	0.52
1:D:1507:ALA:HB2	1:D:1568:VAL:HG11	1.91	0.52
1:D:1825:THR:HG22	1:D:1825:THR:O	2.09	0.52
1:A:362:SER:O	1:A:367:TRP:HB2	2.09	0.52
1:C:1317:ILE:O	1:C:1339:ALA:HB1	2.10	0.52
1:C:1349:ILE:HG12	1:C:1354:GLU:HB3	1.91	0.52
1:C:1225:LEU:O	1:C:1229:VAL:HG23	2.09	0.52
1:D:1769:LYS:HG2	1:D:1772:TRP:CZ2	2.45	0.51
1:C:1154:LEU:HD21	1:C:1218:ILE:HD11	1.93	0.51
1:C:1404:GLU:HB3	1:D:1903:SER:HB2	1.92	0.51
1:C:1233:LYS:HA	1:C:1238:ASP:OD2	2.11	0.51
1:D:1869:VAL:O	1:D:1892:THR:HB	2.11	0.51
1:C:1288:ARG:HG3	1:C:1289:TYR:CD2	2.46	0.51
1:C:1164:LEU:HD22	1:C:1218:ILE:CD1	2.40	0.51
1:A:370:MET:HA	1:A:394:GLN:HG3	1.93	0.51
1:A:317:ILE:O	1:A:339:ALA:HB1	2.10	0.51
1:D:1684:ARG:O	1:D:1688:GLU:HG3	2.11	0.51
1:D:1534:VAL:HG13	1:D:1611:LEU:HD23	1.93	0.50
1:A:246:ASP:HA	1:A:295:GLU:HB3	1.92	0.50
1:A:334:ILE:HD13	1:A:367:TRP:CE2	2.46	0.50
1:C:1164:LEU:HD21	1:C:1169:PHE:CE2	2.47	0.50
1:D:1797:PRO:HD2	1:D:1806:TRP:CH2	2.47	0.50
1:C:1403:SER:HB2	1:D:1904:GLU:HB3	1.91	0.50
1:D:1667:GLN:HG2	1:D:1668:GLU:HG3	1.92	0.50
1:D:1597:ASP:OD2	1:D:1602:LYS:HA	2.12	0.50
1:B:654:LEU:HD23	1:B:654:LEU:C	2.33	0.49
1:A:46:LEU:HD23	1:A:103:SER:HA	1.93	0.49
1:B:901:ALA:O	1:B:902:ARG:HB2	2.11	0.49
1:C:1169:PHE:HB3	1:C:1228:ILE:HD11	1.94	0.49
1:A:401:ALA:O	1:A:402:ARG:HB2	2.12	0.49
1:A:252:PHE:HB3	1:A:262:PHE:CD2	2.48	0.49
1:D:1901:ALA:O	1:D:1902:ARG:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:ASP:OD2	1:A:271:LYS:HG3	2.14	0.48
1:B:849:ILE:HG12	1:B:854:GLU:HB3	1.95	0.48
1:C:1222:GLU:OE2	1:C:1285:LEU:HD21	2.13	0.48
1:C:1264:ASN:HD22	1:C:1267:SER:HA	1.79	0.48
1:B:597:ASP:OD2	1:B:602:LYS:HA	2.13	0.48
1:A:273:LEU:HD22	1:A:277:GLN:HB3	1.95	0.47
1:A:403:SER:OG	1:B:904:GLU:HB3	2.14	0.47
1:D:1815:ILE:HD13	1:D:1815:ILE:N	2.29	0.47
1:D:1597:ASP:HB2	1:D:1610:ILE:HD11	1.95	0.47
1:B:786:MET:HE1	1:B:815:ILE:HG12	1.96	0.47
1:D:1796:ASP:HA	1:D:1806:TRP:CH2	2.50	0.47
1:D:1808:HIS:HE1	4:D:2554:HOH:O	1.98	0.47
1:A:307:SER:O	1:A:311:LYS:HG2	2.15	0.47
1:D:1870:MET:HA	1:D:1894:GLN:HG3	1.96	0.47
1:A:152:ASN:OD1	1:A:168:GLU:HG2	2.15	0.47
1:D:1505:VAL:HG22	1:D:1523:LEU:CD1	2.45	0.47
1:C:1404:GLU:HB3	1:D:1903:SER:CB	2.45	0.47
1:B:546:LEU:HD23	1:B:603:SER:HA	1.97	0.47
1:A:387:VAL:HG21	1:A:395:ILE:HB	1.97	0.47
1:A:259:ASP:OD2	1:A:262:PHE:HA	2.15	0.46
1:C:1247:CYS:SG	1:C:1294:ILE:HD11	2.55	0.46
1:D:1640:LYS:HD3	1:D:1891:ARG:NH1	2.31	0.46
1:B:870:MET:HA	1:B:894:GLN:HG3	1.97	0.46
1:D:1752:PHE:HB3	1:D:1762:PHE:CD2	2.51	0.46
1:B:540:THR:HB	1:B:544:GLU:CD	2.36	0.46
1:A:87:LYS:HG2	4:A:2681:HOH:O	2.15	0.46
1:C:1046:LEU:HD23	1:C:1103:SER:HA	1.97	0.46
1:A:410:ASN:HA	1:A:413:LEU:HD12	1.97	0.45
1:B:705:ALA:HB1	1:B:716:PRO:HB3	1.98	0.45
1:D:1578:LYS:HE3	4:D:3299:HOH:O	2.17	0.45
1:A:225:LEU:O	1:A:229:VAL:HG23	2.17	0.45
1:A:148:VAL:HG22	1:A:425:PHE:CZ	2.52	0.45
1:A:369:VAL:O	1:A:392:THR:HB	2.16	0.45
1:D:1837:LYS:HE3	4:D:2771:HOH:O	2.17	0.44
1:C:1294:ILE:O	1:C:1294:ILE:HG23	2.18	0.44
1:C:1299:ALA:HB1	4:C:2686:HOH:O	2.18	0.44
1:A:215:ALA:HB3	1:B:707:ASN:OD1	2.17	0.44
1:D:1546:LEU:HD23	1:D:1603:SER:HA	2.00	0.44
1:D:1904:GLU:HG2	1:D:1905:ARG:NE	2.33	0.44
1:D:1862:SER:O	1:D:1867:TRP:HB2	2.17	0.44
1:A:326:ASN:O	1:A:330:ILE:HG13	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1373:HIS:CD2	1:C:1373:HIS:H	2.34	0.44
1:C:1326:ASN:HB3	1:C:1329:ARG:CG	2.48	0.44
1:C:1299:ALA:HB3	1:C:1302:ASP:CB	2.48	0.43
1:A:299:ALA:HB3	1:A:302:ASP:CB	2.48	0.43
1:C:1151:LEU:O	1:C:1170:MET:HA	2.17	0.43
4:A:2892:HOH:O	1:B:660:ALA:HA	2.17	0.43
1:B:534:VAL:HG13	1:B:611:LEU:HD23	2.01	0.43
1:C:1007:ALA:HB2	1:C:1068:VAL:HG11	1.99	0.43
1:B:862:SER:O	1:B:867:TRP:HB2	2.17	0.43
1:C:1072:ILE:HD12	1:C:1072:ILE:H	1.83	0.43
1:A:403:SER:CB	1:B:904:GLU:HB3	2.48	0.43
1:D:1656:GLY:HA2	1:D:1660:ALA:HB2	2.01	0.43
1:C:1334:ILE:HD13	1:C:1367:TRP:CE2	2.54	0.42
1:C:1325:THR:O	1:C:1325:THR:HG22	2.19	0.42
1:A:383:ILE:HD11	1:A:408:LYS:HG2	2.01	0.42
1:C:1097:ASP:OD2	1:C:1102:LYS:HA	2.18	0.42
1:C:1313:ALA:HB1	1:C:1317:ILE:HD11	2.00	0.42
1:A:384:ALA:O	1:A:387:VAL:HG12	2.19	0.42
1:A:97:ASP:OD2	1:A:102:LYS:HA	2.20	0.42
1:A:403:SER:HB2	1:B:904:GLU:HB3	2.01	0.42
1:D:1573:ALA:N	1:D:1574:PRO:HD2	2.35	0.42
1:A:300:GLU:O	1:A:322:LEU:HD12	2.20	0.42
1:B:571:VAL:HG11	1:B:605:LEU:HD21	2.02	0.42
1:A:294:ILE:CD1	1:A:297:PRO:HB3	2.50	0.42
1:A:249:SER:HA	1:A:252:PHE:CZ	2.55	0.42
1:B:752:PHE:HB3	1:B:762:PHE:CD2	2.54	0.42
1:A:278:LEU:HG	1:A:282:TYR:CE2	2.55	0.42
1:D:1873:HIS:C	1:D:1873:HIS:CD2	2.93	0.42
1:C:1295:GLU:OE2	1:C:1343:LEU:HD22	2.20	0.41
1:A:218:ILE:CG2	1:A:223:GLU:HB3	2.51	0.41
1:C:1403:SER:CB	1:D:1904:GLU:HB3	2.50	0.41
1:D:1716:PRO:HG2	1:D:1718:ILE:HG13	2.02	0.41
1:C:1296:ASP:OD2	1:C:1320:ASP:HB3	2.20	0.41
1:D:1786:MET:HE1	1:D:1813:ALA:O	2.21	0.41
1:B:796:ASP:HA	1:B:806:TRP:CH2	2.56	0.41
1:A:10:VAL:O	1:A:17:PRO:HA	2.21	0.41
1:A:73:ALA:N	1:A:74:PRO:HD2	2.36	0.41
1:B:694:LYS:HE3	1:B:706:GLY:O	2.21	0.41
1:C:1401:ALA:O	1:C:1402:ARG:CB	2.68	0.41
1:A:299:ALA:HB1	4:A:2900:HOH:O	2.21	0.41
1:A:42:VAL:HG12	1:A:300:GLU:OE1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ARG:HH12	1:A:375:SER:CB	2.34	0.41
1:A:128:PRO:O	1:A:131:LYS:HB3	2.21	0.41
1:D:1856:ILE:O	1:D:1860:GLN:HG3	2.20	0.41
1:C:1252:PHE:HB2	1:C:1259:ASP:O	2.21	0.41
1:B:552:ASP:O	1:B:558:GLY:HA2	2.21	0.41
1:C:1156:GLY:HA3	1:C:1164:LEU:HD23	2.03	0.41
1:D:1817:ILE:O	1:D:1839:ALA:HB1	2.21	0.41
1:A:326:ASN:HB3	1:A:329:ARG:CG	2.52	0.40
1:B:799:ALA:HB1	4:B:3253:HOH:O	2.19	0.40
1:C:1222:GLU:H	1:C:1222:GLU:CD	2.25	0.40
1:B:919:LEU:HD13	1:B:923:ALA:HB2	2.02	0.40
1:B:763:LYS:HB2	4:B:3295:HOH:O	2.22	0.40
1:B:517:PRO:HG2	1:B:556:TRP:CG	2.55	0.40
1:A:373:HIS:CD2	1:A:373:HIS:H	2.38	0.40
1:A:67:ASN:HD22	1:A:67:ASN:HA	1.63	0.40
1:C:1288:ARG:HG3	1:C:1289:TYR:CE2	2.56	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	434/436 (100%)	424 (98%)	9 (2%)	1 (0%)	52	53
1	B	434/436 (100%)	420 (97%)	13 (3%)	1 (0%)	52	53
1	C	434/436 (100%)	419 (96%)	15 (4%)	0	100	100
1	D	434/436 (100%)	421 (97%)	13 (3%)	0	100	100
All	All	1736/1744 (100%)	1684 (97%)	50 (3%)	2 (0%)	56	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	402	ARG
1	B	902	ARG

### 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	343/343 (100%)	337 (98%)	6 (2%)	68	74
1	B	343/343 (100%)	337 (98%)	6 (2%)	68	74
1	C	343/343 (100%)	331 (96%)	12 (4%)	43	44
1	D	343/343 (100%)	339 (99%)	4 (1%)	78	84
All	All	1372/1372 (100%)	1344 (98%)	28 (2%)	63	68

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	143	PRO
1	A	238	ASP
1	A	255	ASP
1	A	347	ASN
1	A	373	HIS
1	B	580	ASN
1	B	626	ASN
1	B	652	ASN
1	B	835	GLU
1	B	847	ASN
1	B	873	HIS
1	C	1080	ASN
1	C	1143	PRO
1	C	1151	LEU
1	C	1152	ASN
1	C	1218	ILE
1	C	1227	LEU
1	C	1255	ASP
1	C	1294	ILE

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Mol	Chain	Res	Type
1	C	1347	ASN
1	C	1373	HIS
1	C	1392	THR
1	C	1428	GLU
1	D	1652	ASN
1	D	1815	ILE
1	D	1847	ASN
1	D	1873	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	A	80	ASN
1	A	191	HIS
1	B	516	ASN
1	B	580	ASN
1	B	626	ASN
1	B	652	ASN
1	B	659	HIS
1	B	808	HIS
1	B	922	ASN
1	C	1067	ASN
1	C	1080	ASN
1	C	1264	ASN
1	C	1308	HIS
1	D	1580	ASN
1	D	1626	ASN
1	D	1652	ASN
1	D	1655	ASN
1	D	1692	ASN
1	D	1808	HIS
1	D	1922	ASN

### 5.3.3 RNA ⓘ

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 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	PAH	A	440	2	8,8,8	2.02	2 (25%)	11,11,11	1.56	3 (27%)
3	PAH	B	940	2	8,8,8	2.20	2 (25%)	11,11,11	1.57	3 (27%)
3	PAH	C	1440	2	8,8,8	2.28	2 (25%)	11,11,11	1.57	3 (27%)
3	PAH	D	1940	2	8,8,8	2.34	2 (25%)	11,11,11	1.45	2 (18%)

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	PAH	A	440	2	-	0/7/7/7	0/0/0/0
3	PAH	B	940	2	-	0/7/7/7	0/0/0/0
3	PAH	C	1440	2	-	0/7/7/7	0/0/0/0
3	PAH	D	1940	2	-	0/7/7/7	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	440	PAH	O2-C2	3.18	1.29	1.23
3	D	1940	PAH	O2-C2	3.42	1.30	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1440	PAH	O2-C2	3.42	1.30	1.23
3	B	940	PAH	O2-C2	3.64	1.30	1.23
3	A	440	PAH	P-C1	4.09	1.85	1.79
3	B	940	PAH	P-C1	4.29	1.86	1.79
3	C	1440	PAH	P-C1	4.84	1.86	1.79
3	D	1940	PAH	P-C1	4.93	1.86	1.79

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1440	PAH	O3-N3-C2	-2.86	115.23	119.56
3	A	440	PAH	O3-N3-C2	-2.79	115.33	119.56
3	B	940	PAH	O3-N3-C2	-2.79	115.34	119.56
3	D	1940	PAH	O3-N3-C2	-2.74	115.42	119.56
3	A	440	PAH	O2-C2-N3	-2.43	120.66	123.53
3	B	940	PAH	O2-C2-N3	-2.41	120.69	123.53
3	D	1940	PAH	O2-C2-N3	-2.24	120.89	123.53
3	C	1440	PAH	O2-C2-N3	-2.23	120.91	123.53
3	A	440	PAH	C1-C2-N3	2.06	118.73	114.59
3	C	1440	PAH	C1-C2-N3	2.15	118.91	114.59
3	B	940	PAH	C1-C2-N3	2.16	118.94	114.59

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 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.