



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2016 – 10:33 AM EDT

PDB ID : 5B3A  
Title : Crystal Structure of O-Phosphoserine Sulfhydrylase from *Aeropyrum pernix* in Complexed with the alpha-Aminoacrylate Intermediate  
Authors : Nakamura, T.; Takeda, E.; Kawai, Y.; Kataoka, M.; Ishikawa, K.  
Deposited on : 2016-02-12  
Resolution : 2.14 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.

---

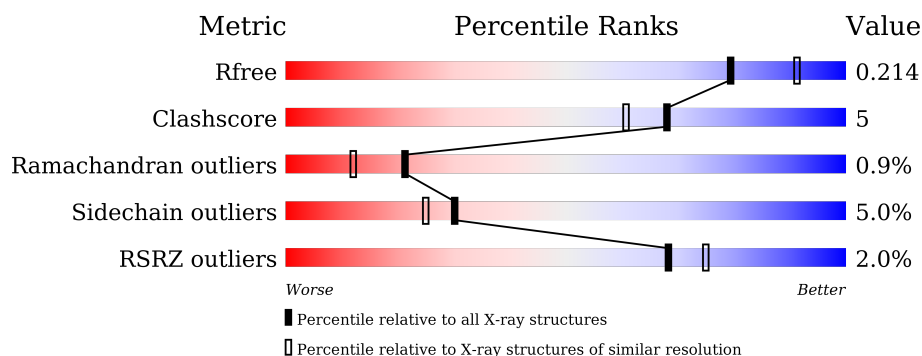
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027939  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027939

**i**

## X-RAY DIFFRACTION

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))
$R_{free}$	91344	1693 (2.16-2.12)
Clashscore	102246	1824 (2.16-2.12)
Ramachandran outliers	100387	1798 (2.16-2.12)
Sidechain outliers	100360	1798 (2.16-2.12)
RSRZ outliers	91569	1699 (2.16-2.12)

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.

Mol	Chain	Length	Quality of chain
1	A	389	<div> <div> <div>3%</div> <div>85%</div> <div>12%</div> <div>...</div> </div> </div>
1	B	389	<div> <div> <div>%</div> <div>85%</div> <div>11%</div> <div>...</div> </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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MPD	B	401	-	-	-	X

## 2 Entry composition [i](#)

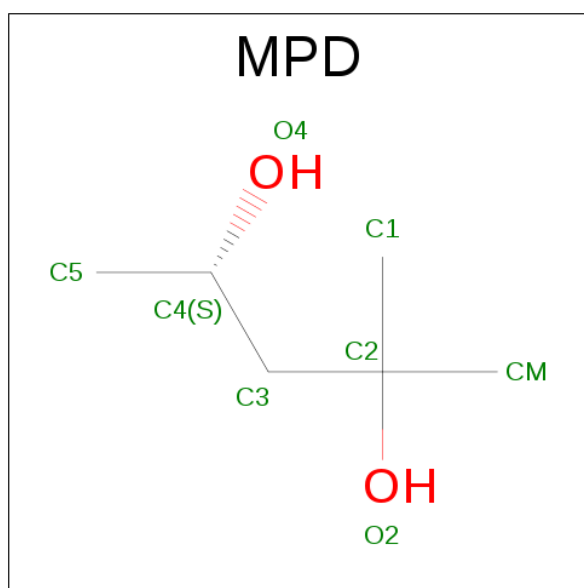
There are 4 unique types of molecules in this entry. The entry contains 6326 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Protein CysO.

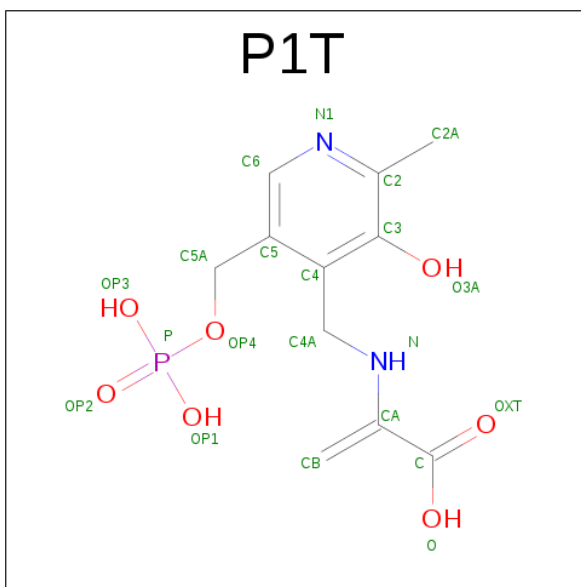
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	0	4	0
			2945	1868	518	550	9			
1	B	383	Total	C	N	O	S	0	9	0
			2975	1890	518	558	9			

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula:  $C_6H_{14}O_2$ ).



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

- Molecule 3 is 2-[(3-HYDROXY-2-METHYL-5-[(PHOSPHONOOXY)METHYL]PYRIDIN-4-YL)METHYL]AMINO]ACRYLIC ACID (three-letter code: P1T) (formula:  $C_{11}H_{15}N_2O_7P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			21	11	2	7	1		
3	B	1	Total	C	N	O	P	0	0
			21	11	2	7	1		

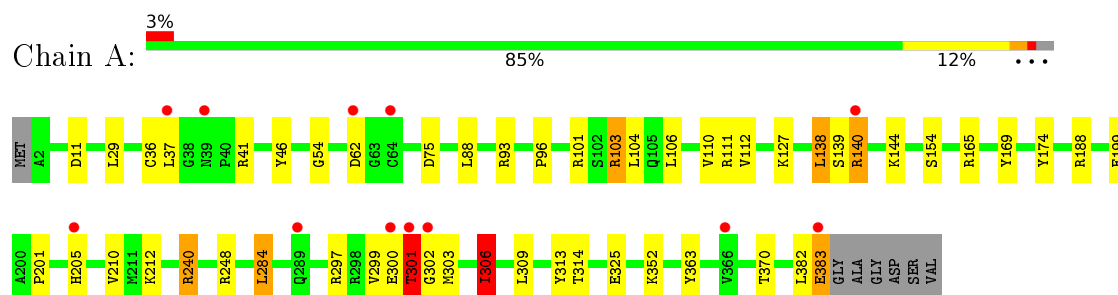
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	162	Total	O	0	0
			162	162		
4	B	186	Total	O	0	0
			186	186		

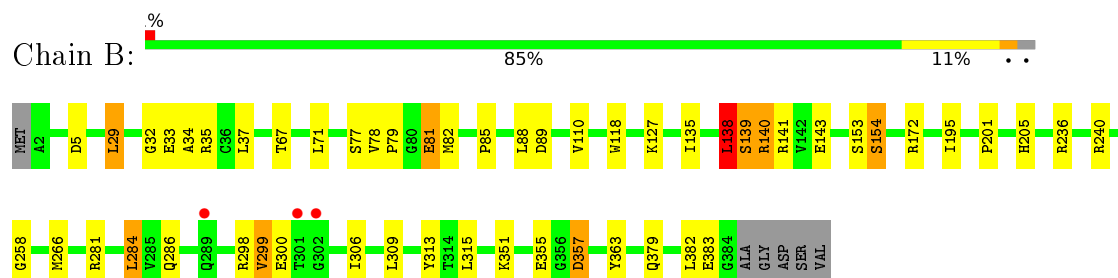
### 3 Residue-property plots [i](#)

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.

#### • Molecule 1: Protein CysO



#### • Molecule 1: Protein CysO



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.15Å 74.15Å 275.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.92 – 2.14 32.92 – 2.14	Depositor EDS
% Data completeness (in resolution range)	98.0 (32.92-2.14) 98.0 (32.92-2.14)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	18.63 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.142 , 0.203 0.154 , 0.214	Depositor DCC
$R_{free}$ test set	2161 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.2	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 48.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6326	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MPD, P1T

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	0.90	0/3010	1.00	13/4084 (0.3%)
1	B	0.92	2/3059 (0.1%)	0.98	7/4150 (0.2%)
All	All	0.91	2/6069 (0.0%)	0.99	20/8234 (0.2%)

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	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	81	GLU	CG-CD	6.47	1.61	1.51
1	B	143	GLU	CG-CD	5.79	1.60	1.51

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	297	ARG	NE-CZ-NH2	8.18	124.39	120.30
1	A	297	ARG	NE-CZ-NH1	-7.43	116.58	120.30
1	A	188	ARG	NE-CZ-NH2	7.34	123.97	120.30
1	A	93	ARG	NE-CZ-NH1	7.12	123.86	120.30
1	B	298	ARG	NE-CZ-NH1	-6.84	116.88	120.30
1	A	93	ARG	NE-CZ-NH2	-6.64	116.98	120.30
1	A	111	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	A	103	ARG	NE-CZ-NH2	-6.42	117.09	120.30
1	A	103	ARG	NE-CZ-NH1	6.30	123.45	120.30

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	ILE	CB-CA-C	-5.99	99.61	111.60
1	B	139	SER	N-CA-C	5.98	127.14	111.00
1	B	139	SER	CA-C-N	5.92	130.22	117.20
1	B	139	SER	C-N-CA	5.60	135.69	121.70
1	A	101	ARG	NE-CZ-NH1	-5.49	117.56	120.30
1	B	298	ARG	CG-CD-NE	-5.30	100.67	111.80
1	B	138	LEU	CA-CB-CG	5.30	127.48	115.30
1	A	75	ASP	CB-CG-OD2	-5.28	113.55	118.30
1	B	29	LEU	CB-CG-CD1	5.17	119.78	111.00
1	A	188	ARG	NE-CZ-NH1	-5.17	117.72	120.30
1	A	240	ARG	NE-CZ-NH1	5.15	122.87	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	138	LEU	Peptide
1	A	382	LEU	Peptide

## 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	2945	0	2965	25	0
1	B	2975	0	2999	28	0
2	A	8	0	14	2	0
2	B	8	0	14	3	0
3	A	21	0	11	1	0
3	B	21	0	12	1	0
4	A	162	0	0	5	2
4	B	186	0	0	6	3
All	All	6326	0	6015	56	3

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 (56) 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:A:300:GLU:O	1:A:301:THR:HB	1.43	1.09
1:B:236:ARG:HD2	4:B:624:HOH:O	1.69	0.92
1:A:300:GLU:O	1:A:301:THR:CB	2.23	0.86
1:B:382:LEU:O	4:B:501:HOH:O	1.97	0.82
2:B:401:MPD:H12	2:B:401:MPD:H52	1.61	0.81
1:B:5[A]:ASP:OD1	4:B:502:HOH:O	1.99	0.79
1:A:37:LEU:HD21	1:A:46:TYR:CD2	2.22	0.74
1:A:139:SER:HA	1:A:169:TYR:OH	1.97	0.64
1:A:96:PRO:O	2:A:401:MPD:HM2	1.97	0.64
1:B:33[B]:GLU:OE1	4:B:504:HOH:O	2.15	0.64
1:B:33[B]:GLU:OE2	1:B:67:THR:HA	2.05	0.57
1:A:299:VAL:O	1:A:301:THR:HG22	2.06	0.56
1:B:138:LEU:HD23	1:B:138:LEU:C	2.26	0.56
1:B:284:LEU:HD22	1:B:313:TYR:HB2	1.90	0.53
1:A:301:THR:HG23	1:A:302:GLY:N	2.24	0.52
1:B:138:LEU:C	1:B:140:ARG:HB3	2.29	0.52
1:A:201:PRO:HD2	1:A:205:HIS:CG	2.45	0.51
1:A:303:MET:HB2	1:A:306:ILE:HD11	1.93	0.51
1:A:284:LEU:HD22	1:A:313:TYR:HB2	1.93	0.50
1:A:240:ARG:HD2	4:A:521:HOH:O	2.13	0.49
2:B:401:MPD:H52	2:B:401:MPD:C1	2.38	0.48
1:A:240:ARG:CD	4:A:521:HOH:O	2.63	0.47
1:A:248[A]:ARG:NH1	1:B:81:GLU:OE2	2.48	0.47
1:B:172:ARG:HD3	1:B:195:ILE:HD12	1.97	0.46
1:A:140:ARG:HB2	1:A:140:ARG:NH1	2.31	0.46
1:B:67:THR:O	1:B:71:LEU:HG	2.16	0.45
1:A:174:TYR:CD1	1:A:210:VAL:HG22	2.51	0.45
1:A:127:LYS:HD3	3:A:402:P1T:H4A2	1.99	0.45
1:B:286:GLN:HG2	1:B:315:LEU:HD22	2.00	0.44
1:B:201:PRO:HD2	1:B:205:HIS:CG	2.53	0.44
1:B:110:VAL:HG13	1:B:363:TYR:CD1	2.53	0.44
1:B:35:ARG:NH1	4:B:515:HOH:O	2.51	0.44
1:B:78:VAL:HG22	1:B:82:MET:SD	2.58	0.43
1:B:85:PRO:HD2	1:B:89:ASP:OD1	2.17	0.43
1:B:139:SER:N	1:B:140:ARG:CB	2.81	0.43
1:B:258:GLY:HA3	1:B:266:MET:SD	2.58	0.43
1:A:165:ARG:HD3	4:A:590:HOH:O	2.19	0.42
1:B:37:LEU:HA	1:B:37:LEU:HD23	1.81	0.42
1:A:383:GLU:HA	1:A:383:GLU:OE2	2.19	0.42
1:A:54:GLY:HA3	1:A:240:ARG:HD3	2.02	0.42
2:B:401:MPD:H4	4:B:515:HOH:O	2.19	0.42
1:B:32:GLY:O	1:B:33[B]:GLU:OE2	2.38	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:LYS:HD3	3:B:402:P1T:H4A2	2.02	0.41
1:B:34:ALA:HB3	1:B:71:LEU:HD11	2.00	0.41
1:A:240:ARG:NE	4:A:521:HOH:O	2.53	0.41
1:B:351:LYS:O	1:B:355:GLU:HG3	2.20	0.41
1:A:314:THR:HG22	1:A:352:LYS:HE3	2.03	0.41
1:B:153:SER:O	1:B:154:SER:CB	2.69	0.41
1:A:110:VAL:HG13	1:A:363:TYR:CD1	2.56	0.41
1:A:106:LEU:HD13	1:A:112:VAL:CG2	2.51	0.40
2:A:401:MPD:HM3	4:A:535:HOH:O	2.22	0.40
1:B:135:ILE:O	1:B:138:LEU:HD22	2.21	0.40
1:A:11:ASP:CG	1:A:41[A]:ARG:HH22	2.23	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:550:HOH:O	4:B:680:HOH:O[4_555]	1.63	0.57
4:B:568:HOH:O	4:B:672:HOH:O[5_654]	1.79	0.41
4:A:600:HOH:O	4:B:515:HOH:O[5_654]	2.18	0.02

## 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	384/389 (99%)	374 (97%)	8 (2%)	2 (0%)	34	26
1	B	390/389 (100%)	371 (95%)	13 (3%)	6 (2%)	13	5
All	All	774/778 (100%)	745 (96%)	21 (3%)	8 (1%)	21	10

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	SER
1	A	301	THR
1	B	141	ARG
1	B	154	SER
1	B	300[A]	GLU
1	B	300[B]	GLU
1	B	140	ARG
1	B	299	VAL

### 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	313/313 (100%)	296 (95%)	17 (5%)	27	21
1	B	318/313 (102%)	302 (95%)	16 (5%)	30	25
All	All	631/626 (101%)	598 (95%)	33 (5%)	30	23

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LEU
1	A	36	CYS
1	A	62	ASP
1	A	88	LEU
1	A	104	LEU
1	A	138	LEU
1	A	140	ARG
1	A	144	LYS
1	A	199	GLU
1	A	212	LYS
1	A	284	LEU
1	A	301	THR
1	A	306	ILE
1	A	309	LEU
1	A	325	GLU
1	A	370	THR

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	383	GLU
1	B	29	LEU
1	B	77	SER
1	B	79	PRO
1	B	88	LEU
1	B	118	TRP
1	B	138	LEU
1	B	240	ARG
1	B	281[A]	ARG
1	B	281[B]	ARG
1	B	284	LEU
1	B	306	ILE
1	B	309	LEU
1	B	357[A]	ASP
1	B	357[B]	ASP
1	B	379	GLN
1	B	383	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS

### 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](#)

4 ligands are modelled in this entry.

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	$\# Z  > 2$
2	MPD	A	401	-	6,7,7	0.71	0	6,10,10	1.04	0
3	P1T	A	402	-	19,21,21	1.57	4 (21%)	23,30,30	1.82	7 (30%)
2	MPD	B	401	-	6,7,7	0.77	0	6,10,10	0.82	0
3	P1T	B	402	-	19,21,21	1.97	5 (26%)	23,30,30	2.05	4 (17%)

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
2	MPD	A	401	-	-	0/5/5/5	0/0/0/0
3	P1T	A	402	-	-	0/10/15/15	0/1/1/1
2	MPD	B	401	-	-	0/5/5/5	0/0/0/0
3	P1T	B	402	-	-	0/10/15/15	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	P1T	C3-C2	-5.63	1.36	1.40
3	A	402	P1T	C4A-N	-3.16	1.39	1.46
3	B	402	P1T	C4A-N	-2.91	1.40	1.46
3	B	402	P1T	C4A-C4	-2.55	1.49	1.51
3	A	402	P1T	C-CA	-2.24	1.48	1.52
3	A	402	P1T	C2A-C2	2.17	1.54	1.50
3	B	402	P1T	P-OP4	2.54	1.67	1.59
3	B	402	P1T	CB-CA	2.94	1.39	1.33
3	A	402	P1T	CB-CA	3.33	1.39	1.33

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	P1T	C4A-C4-C5	-5.70	114.26	119.72
3	A	402	P1T	C4A-C4-C5	-5.32	114.61	119.72

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	P1T	C5-C6-N1	-2.80	118.96	123.86
3	A	402	P1T	OP3-P-OP1	-2.37	98.74	107.44
3	A	402	P1T	C3-C2-N1	-2.14	118.01	120.69
3	A	402	P1T	C6-N1-C2	2.02	123.32	119.26
3	A	402	P1T	C2A-C2-C3	2.27	123.20	120.90
3	B	402	P1T	C4-C4A-N	2.41	116.42	111.04
3	A	402	P1T	OP3-P-OP4	2.55	114.16	106.72
3	A	402	P1T	O3A-C3-C2	2.73	121.53	117.53
3	B	402	P1T	O3A-C3-C2	4.94	124.77	117.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	MPD	2	0
3	A	402	P1T	1	0
2	B	401	MPD	3	0
3	B	402	P1T	1	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	382/389 (98%)	-0.20	12 (3%) 52 62	9, 16, 38, 77	0
1	B	383/389 (98%)	-0.25	3 (0%) 87 90	9, 16, 33, 62	0
All	All	765/778 (98%)	-0.22	15 (1%) 68 75	9, 16, 36, 77	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	302	GLY	7.7
1	A	301	THR	5.7
1	B	301	THR	4.8
1	A	62	ASP	4.3
1	A	64	CYS	2.9
1	A	140	ARG	2.8
1	B	302	GLY	2.7
1	A	289	GLN	2.6
1	A	37	LEU	2.6
1	A	300	GLU	2.5
1	A	383	GLU	2.4
1	A	39	ASN	2.2
1	A	205	HIS	2.1
1	A	366	VAL	2.1
1	B	289	GLN	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	MPD	B	401	8/8	0.85	0.19	3.78	41,46,48,52	0
2	MPD	A	401	8/8	0.93	0.14	2.00	13,16,16,20	0
3	P1T	A	402	21/21	0.99	0.10	0.02	10,12,13,13	0
3	P1T	B	402	21/21	0.98	0.11	-0.02	10,11,12,12	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.