



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 10:44 PM GMT

PDB ID : 3Q3M  
Title : Toluene 4 monooxygenase HD Complex with Inhibitor 4-Bromobenzoate  
Authors : Acheson, J.F.; Bailey, L.J.; Fox, B.G.  
Deposited on : 2010-12-22  
Resolution : 1.75 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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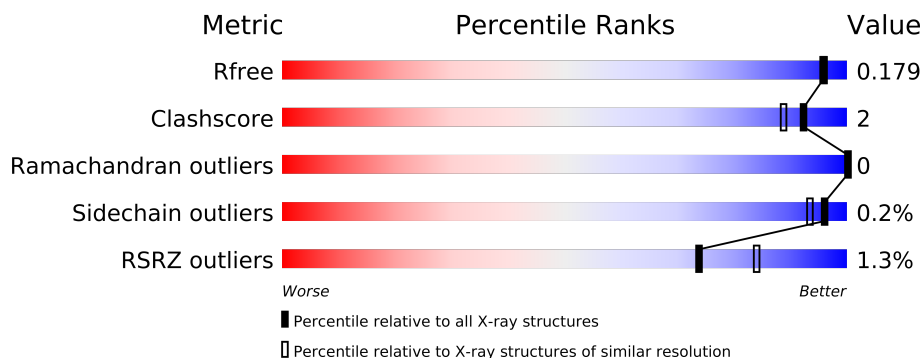
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.75 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))
$R_{free}$	66092	1657 (1.76-1.72)
Clashscore	79885	1881 (1.76-1.72)
Ramachandran outliers	78287	1859 (1.76-1.72)
Sidechain outliers	78261	1859 (1.76-1.72)
RSRZ outliers	66119	1658 (1.76-1.72)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	500	
1	D	500	
2	B	307	
2	F	307	
3	C	84	
3	G	84	
4	E	103	
4	H	103	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
6	Z82	A	510	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
6	Z82	D	511	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 18516 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Toluene-4-monooxygenasesystem protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	491	Total	C	N	O	S	0	4	0
			4065	2603	687	751	24			
1	D	491	Total	C	N	O	S	0	12	0
			4127	2643	697	762	25			

- Molecule 2 is a protein called Toluene-4-monooxygenasesystem protein E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	305	Total	C	N	O	S	0	7	0
			2578	1629	447	487	15			
2	F	305	Total	C	N	O	S	0	9	0
			2601	1643	452	490	16			

- Molecule 3 is a protein called Toluene-4-monooxygenasesystem protein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	82	Total	C	N	O	S	0	0	0
			654	412	117	121	4			
3	G	82	Total	C	N	O	S	0	1	0
			662	416	118	124	4			

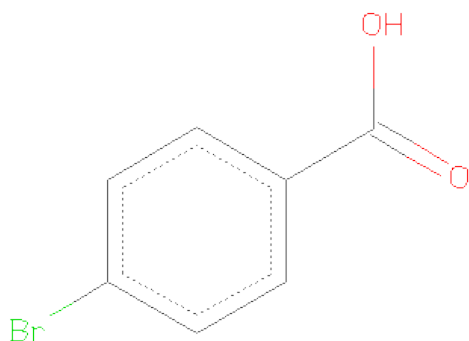
- Molecule 4 is a protein called Toluene-4-monooxygenasesystem protein D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	102	Total	C	N	O	S	0	1	0
			815	508	141	163	3			
4	H	102	Total	C	N	O	S	0	1	0
			815	508	141	163	3			

- Molecule 5 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Fe	0	0
			2	2		
5	D	2	Total	Fe	0	0
			2	2		

- Molecule 6 is 4-BROMOBENZOIC ACID (three-letter code: Z82) (formula: C<sub>7</sub>H<sub>5</sub>BrO<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	Br	C	O	0	0
			10	1	7	2		
6	B	1	Total	Br	C	O	0	0
			10	1	7	2		
6	B	1	Total	Br	C	O	0	0
			10	1	7	2		
6	E	1	Total	Br	C	O	0	0
			10	1	7	2		
6	D	1	Total	Br	C	O	0	0
			10	1	7	2		
6	F	1	Total	Br	C	O	0	0
			10	1	7	2		
6	H	1	Total	Br	C	O	0	0
			10	1	7	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	487	Total	O	0	0
			487	487		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	302	Total 302	O 302	0	0
7	C	74	Total 74	O 74	0	0
7	E	148	Total 148	O 148	0	0
7	D	531	Total 531	O 531	0	0
7	F	357	Total 357	O 357	0	0
7	G	85	Total 85	O 85	0	0
7	H	141	Total 141	O 141	0	0

### 3 Residue-property plots

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.

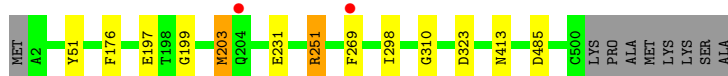
- Molecule 1: Toluene-4-monooxygenasesystem protein A

Chain A: 



- Molecule 1: Toluene-4-monooxygenasesystem protein A

Chain D: 



- Molecule 2: Toluene-4-monooxygenasesystem protein E

Chain B: 



- Molecule 2: Toluene-4-monooxygenasesystem protein E

Chain F: 



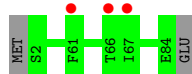
- Molecule 3: Toluene-4-monooxygenasesystem protein B

Chain C: 



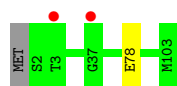
- Molecule 3: Toluene-4-monooxygenasesystem protein B

Chain G: 



- Molecule 4: Toluene-4-monooxygenasesystem protein D

Chain E: 



- Molecule 4: Toluene-4-monooxygenasesystem protein D

Chain H: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.04Å 115.72Å 181.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.09 – 1.75 48.77 – 1.75	Depositor EDS
% Data completeness (in resolution range)	93.6 (34.09-1.75) 97.6 (48.77-1.75)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.35 (at 1.75Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, $R_{free}$	0.143 , 0.178 0.146 , 0.179	Depositor DCC
$R_{free}$ test set	10387 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	14.9	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 39.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 207453 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	18516	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 57.56 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.3267e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: Z82, 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	0.55	0/4194	0.61	1/5694 (0.0%)
1	D	0.55	0/4256	0.62	2/5778 (0.0%)
2	B	0.47	0/2654	0.56	0/3608
2	F	0.53	1/2676 (0.0%)	0.61	0/3634
3	C	0.45	0/666	0.60	0/902
3	G	0.45	0/674	0.56	0/913
4	E	0.53	0/824	0.65	0/1112
4	H	0.49	0/824	0.62	0/1112
All	All	0.52	1/16768 (0.0%)	0.61	3/22753 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	9	MET	C-N	-5.25	1.22	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	203	MET	CG-SD-CE	6.93	111.28	100.20
1	D	251	ARG	NE-CZ-NH1	-5.63	117.48	120.30
1	A	203	MET	CG-SD-CE	5.17	108.47	100.20

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4065	0	0	10	0
1	D	4127	0	4	8	0
2	B	2578	0	9	1	0
2	F	2601	0	0	6	0
3	C	654	0	0	0	0
3	G	662	0	0	0	0
4	E	815	0	0	1	0
4	H	815	0	4	2	0
5	A	2	0	0	0	0
5	D	2	0	0	0	0
6	A	10	0	4	1	0
6	B	20	0	8	0	0
6	D	10	0	4	1	0
6	E	10	0	4	1	0
6	F	10	0	4	1	0
6	H	10	0	4	1	0
7	A	487	0	0	3	0
7	B	302	0	0	1	0
7	C	74	0	0	0	0
7	D	531	0	0	5	0
7	E	148	0	0	1	0
7	F	357	0	0	3	0
7	G	85	0	0	0	0
7	H	141	0	0	1	0
All	All	18516	0	45	29	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

All (29) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
7:D:2014:HOH:O	6:H:104:Z82:BR4	2.21	1.11
1:A:203:MET:CB	1:A:269:PHE:CZ	2.65	0.79
1:D:203:MET:CB	1:D:269:PHE:CZ	2.67	0.77
1:D:413[B]:ASN:ND2	7:D:1577:HOH:O	2.18	0.76
1:A:209:ALA:N	7:A:2027:HOH:O	2.29	0.64
4:H:78:GLU:OE2	7:H:1438:HOH:O	2.15	0.63
2:F:10:ARG:NH1	2:F:22:PRO:CD	2.61	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:251:ARG:NH1	7:D:1753:HOH:O	2.33	0.62
2:F:70:ASP:OD1	7:F:329:HOH:O	2.15	0.60
1:A:80:LYS:NZ	4:H:46:ARG:NH1	2.49	0.60
1:D:199:GLY:O	1:D:298:ILE:CD1	2.51	0.59
7:D:1653:HOH:O	2:F:48:ASP:CB	2.53	0.57
1:A:205:PHE:C	7:A:2027:HOH:O	2.43	0.57
1:D:197:GLU:CD	1:D:231:GLU:OE1	2.44	0.56
1:A:80:LYS:NZ	7:A:1827:HOH:O	2.42	0.52
2:F:104[A]:ARG:NE	7:F:1711:HOH:O	2.43	0.51
1:A:197:GLU:CD	1:A:231:GLU:OE1	2.51	0.49
4:E:78:GLU:OE2	7:E:1709:HOH:O	2.20	0.46
1:D:485:ASP:CB	7:D:2023:HOH:O	2.65	0.45
1:A:251:ARG:NH1	1:A:310:GLY:O	2.50	0.44
1:D:251:ARG:NH1	1:D:310:GLY:O	2.50	0.44
1:A:203:MET:CG	1:A:269:PHE:CZ	3.01	0.43
2:F:287:GLU:OE1	2:F:290:ARG:NH1	2.51	0.43
6:F:328:Z82:H6	7:F:2054:HOH:O	2.19	0.43
2:B:70:ASP:OD1	7:B:334:HOH:O	2.22	0.42
1:D:176:PHE:CE2	6:D:511:Z82:BR4	3.28	0.41
1:A:6:ARG:NH2	6:E:104:Z82:C7	2.84	0.41
1:A:176:PHE:CE2	6:A:510:Z82:BR4	3.29	0.41
2:F:76:ASP:OD1	2:F:263:ARG:NH2	2.54	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	493/500 (99%)	478 (97%)	15 (3%)	0	100	100
1	D	501/500 (100%)	487 (97%)	14 (3%)	0	100	100
2	B	310/307 (101%)	307 (99%)	3 (1%)	0	100	100
2	F	312/307 (102%)	309 (99%)	3 (1%)	0	100	100
3	C	80/84 (95%)	77 (96%)	3 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	G	81/84 (96%)	77 (95%)	4 (5%)	0	100	100
4	E	100/103 (97%)	99 (99%)	1 (1%)	0	100	100
4	H	100/103 (97%)	99 (99%)	1 (1%)	0	100	100
All	All	1977/1988 (99%)	1933 (98%)	44 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	420/423 (99%)	418 (100%)	2 (0%)	94	89
1	D	428/423 (101%)	426 (100%)	2 (0%)	94	89
2	B	283/278 (102%)	283 (100%)	0	100	100
2	F	285/278 (102%)	285 (100%)	0	100	100
3	C	73/75 (97%)	73 (100%)	0	100	100
3	G	74/75 (99%)	74 (100%)	0	100	100
4	E	87/87 (100%)	87 (100%)	0	100	100
4	H	87/87 (100%)	87 (100%)	0	100	100
All	All	1737/1726 (101%)	1733 (100%)	4 (0%)	96	93

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	TYR
1	A	323	ASP
1	D	51	TYR
1	D	323	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains 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 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	Z82	A	510	5	10,10,10	1.83	3 (30%)	13,13,13	1.12	0
6	Z82	B	328	-	10,10,10	1.03	1 (10%)	13,13,13	0.81	0
6	Z82	B	329	-	10,10,10	0.73	0	13,13,13	0.96	0
6	Z82	D	511	5	10,10,10	1.78	2 (20%)	13,13,13	1.87	4 (30%)
6	Z82	E	104	-	10,10,10	1.08	1 (10%)	13,13,13	1.11	0
6	Z82	F	328	-	10,10,10	0.79	0	13,13,13	1.08	1 (7%)
6	Z82	H	104	-	10,10,10	0.79	0	13,13,13	1.16	1 (7%)

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	Z82	A	510	5	-	0/4/4/4	0/1/1/1
6	Z82	B	328	-	-	0/4/4/4	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	Z82	B	329	-	-	0/4/4/4	0/1/1/1
6	Z82	D	511	5	-	0/4/4/4	0/1/1/1
6	Z82	E	104	-	-	0/4/4/4	0/1/1/1
6	Z82	F	328	-	-	0/4/4/4	0/1/1/1
6	Z82	H	104	-	-	0/4/4/4	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	511	Z82	O2-C7	-3.72	1.18	1.30
6	A	510	Z82	O2-C7	-3.26	1.20	1.30
6	A	510	Z82	C6-C1	-3.07	1.33	1.39
6	E	104	Z82	BR4-C4	2.66	1.96	1.90
6	D	511	Z82	C6-C1	-2.35	1.35	1.39
6	B	328	Z82	BR4-C4	2.30	1.95	1.90
6	A	510	Z82	O1-C7	-2.07	1.16	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	511	Z82	C5-C6-C1	-3.09	117.07	120.76
6	D	511	Z82	O2-C7-C1	2.85	122.56	115.01
6	D	511	Z82	C2-C1-C7	-2.84	115.22	120.44
6	D	511	Z82	C6-C1-C2	2.83	122.68	118.63
6	H	104	Z82	C3-C2-C1	-2.25	118.07	120.76
6	F	328	Z82	O2-C7-C1	2.18	120.78	115.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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, 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	491/500 (98%)	-0.32	2 (0%) 90 94	7, 13, 26, 40	1 (0%)
1	D	491/500 (98%)	-0.32	2 (0%) 90 94	6, 12, 25, 39	2 (0%)
2	B	305/307 (99%)	-0.19	7 (2%) 57 67	8, 16, 34, 100	2 (0%)
2	F	305/307 (99%)	-0.23	5 (1%) 68 79	7, 13, 28, 104	0
3	C	82/84 (97%)	0.21	5 (6%) 21 25	11, 22, 43, 56	0
3	G	82/84 (97%)	0.06	3 (3%) 39 47	10, 20, 35, 48	1 (1%)
4	E	102/103 (99%)	-0.34	2 (1%) 62 72	7, 13, 28, 43	0
4	H	102/103 (99%)	-0.41	0 100 100	8, 14, 30, 41	0
All	All	1960/1988 (98%)	-0.25	26 (1%) 74 84	6, 14, 30, 104	6 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	327	ASP	14.8
2	B	327	ASP	13.1
2	B	326	SER	5.2
2	F	326	SER	4.8
2	B	10	ARG	4.5
2	F	305	SER	3.3
2	F	240	TRP	3.2
3	C	2	SER	3.1
2	B	240	TRP	2.7
3	G	61	PHE	2.6
4	E	37	GLY	2.6
1	D	269	PHE	2.6
3	C	72	LEU	2.6
3	C	64	ASP	2.4
2	B	179	ASP	2.3
3	C	67	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	113	GLY	2.2
1	A	269	PHE	2.2
2	B	112	GLU	2.2
1	D	204[A]	GLN	2.1
3	G	67	ILE	2.1
4	E	3	THR	2.1
3	G	66	THR	2.0
1	A	204[A]	GLN	2.0
2	F	188	GLU	2.0
3	C	49	VAL	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
6	Z82	A	510	10/10	0.15	2.70	10,16,20,20	9
6	Z82	D	511	10/10	0.14	2.08	7,20,22,25	8
6	Z82	F	328	10/10	0.15	1.62	27,28,32,47	10
6	Z82	E	104	10/10	0.10	1.53	13,14,20,20	10
6	Z82	B	329	10/10	0.14	1.20	24,27,30,46	10
6	Z82	B	328	10/10	0.14	0.35	17,20,24,32	10
6	Z82	H	104	10/10	0.07	-0.04	10,13,19,19	10
5	FE	A	870	1/1	0.04	-2.66	12,12,12,12	0
5	FE	A	509	1/1	0.03	-2.87	14,14,14,14	0
5	FE	D	510	1/1	0.03	-3.27	13,13,13,13	0
5	FE	D	509	1/1	0.03	-5.39	10,10,10,10	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.