



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 05:47 am GMT

PDB ID : 4S13
Title : Ferulic Acid Decarboxylase (FDC1)
Authors : Lee, S.G.; Bhuiya, M.W.; Yu, O.; Jez, J.M.
Deposited on : 2015-01-07
Resolution : 2.35 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

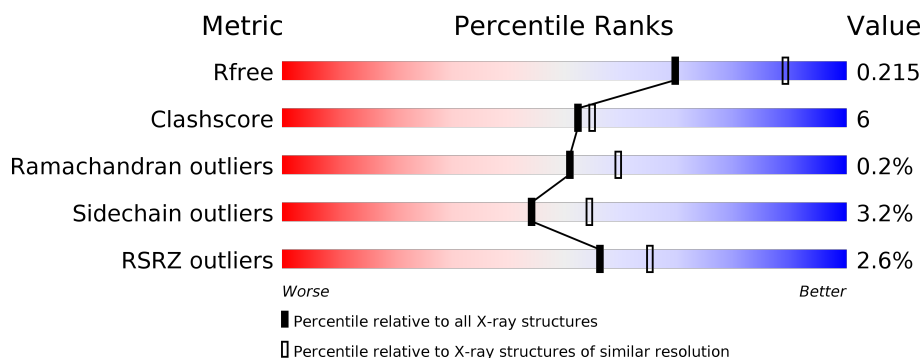
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	100719	1570 (2.36-2.32)
Clashscore	112137	1673 (2.36-2.32)
Ramachandran outliers	110173	1654 (2.36-2.32)
Sidechain outliers	110143	1655 (2.36-2.32)
RSRZ outliers	101464	1576 (2.36-2.32)

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 ≥ 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	503	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>6%</div> </div> </div>
1	B	503	<div> <div>3%</div> <div> <div></div> <div>82%</div> <div>12%</div> <div>5%</div> </div> </div>
1	C	503	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>14%</div> <div>.</div> </div> </div>
1	D	503	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div>5%</div> </div> </div>
1	E	503	<div> <div></div> <div> <div></div> <div>85%</div> <div>12%</div> <div>..</div> </div> </div>
1	F	503	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>17%</div> <div>.</div> </div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	G	503	
1	H	503	

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	4VP	A	601	-	-	-	X
2	4VP	F	601	-	-	-	X
2	4VP	G	601	-	-	-	X

2 Entry composition

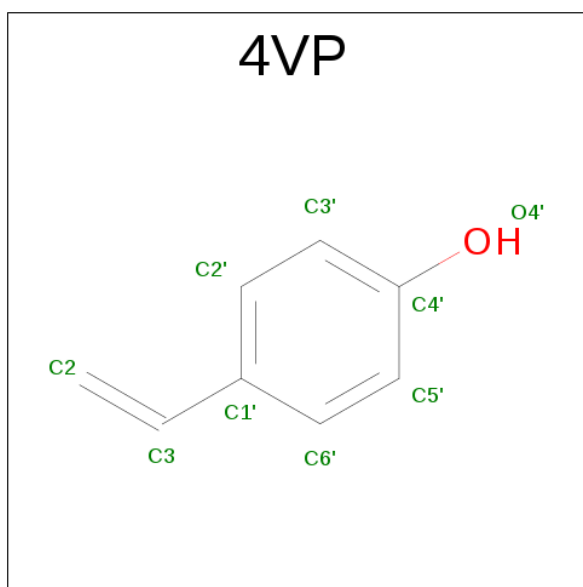
There are 3 unique types of molecules in this entry. The entry contains 31691 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ferulic acid decarboxylase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	S	0	0	0
			3705	2387	600	696	22			
1	B	476	Total	C	N	O	S	0	2	0
			3748	2420	604	701	23			
1	C	501	Total	C	N	O	S	0	0	0
			3933	2536	642	733	22			
1	D	480	Total	C	N	O	S	0	1	0
			3775	2436	610	707	22			
1	E	492	Total	C	N	O	S	0	0	0
			3868	2495	630	721	22			
1	F	501	Total	C	N	O	S	0	0	0
			3933	2536	642	733	22			
1	G	472	Total	C	N	O	S	0	1	0
			3718	2399	600	697	22			
1	H	481	Total	C	N	O	S	0	1	0
			3777	2435	611	709	22			

- Molecule 2 is 4-ETHENYLPHENOL (three-letter code: 4VP) (formula: C₈H₈O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			9	8	1		
2	C	1	Total	C	O	0	0
			9	8	1		
2	E	1	Total	C	O	0	0
			9	8	1		
2	F	1	Total	C	O	0	0
			9	8	1		
2	G	1	Total	C	O	0	0
			9	8	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	152	Total	O	0	0
			152	152		
3	B	190	Total	O	0	0
			190	190		
3	C	131	Total	O	0	0
			131	131		
3	D	185	Total	O	0	0
			185	185		
3	E	165	Total	O	0	0
			165	165		
3	F	137	Total	O	0	0
			137	137		
3	G	100	Total	O	0	0
			100	100		

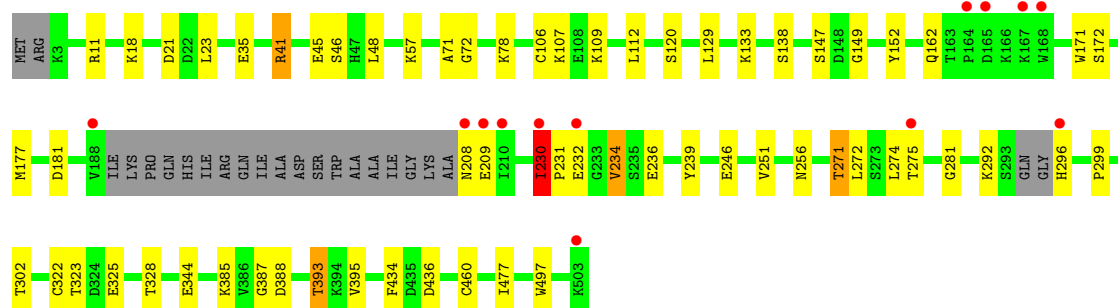
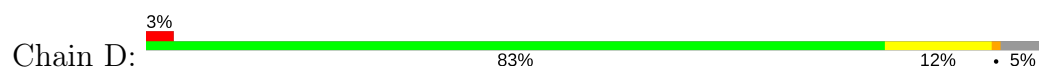
Continued on next page...

Continued from previous page...

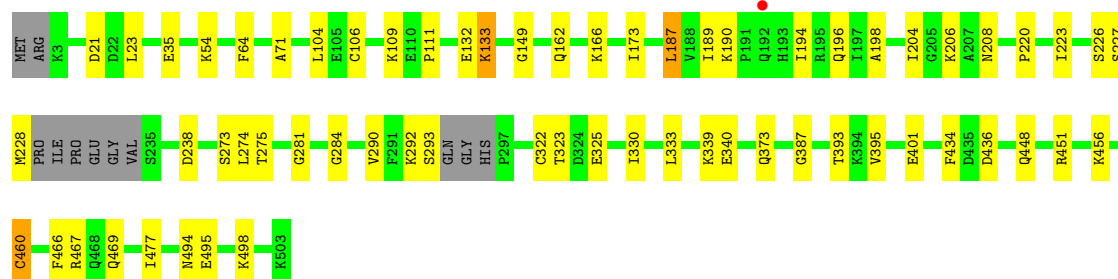
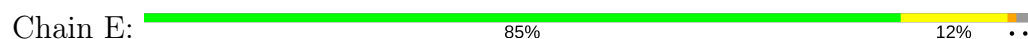
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	129	Total 129	O 129	0	0



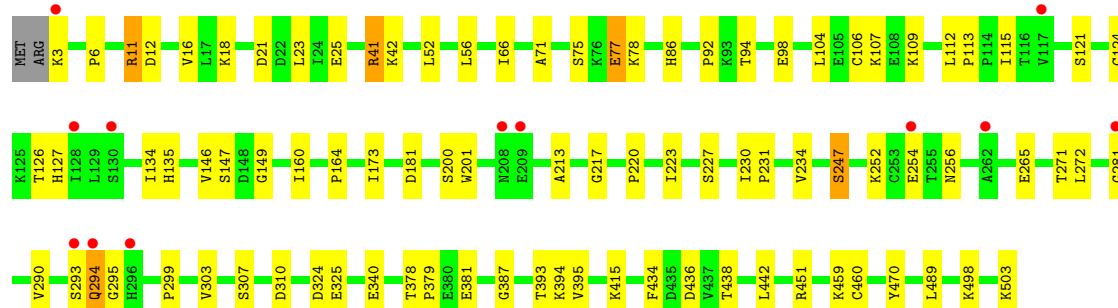
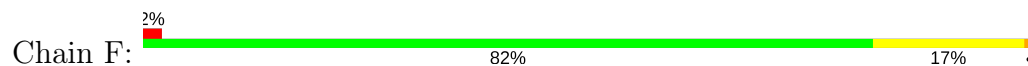
• Molecule 1: Ferulic acid decarboxylase 1



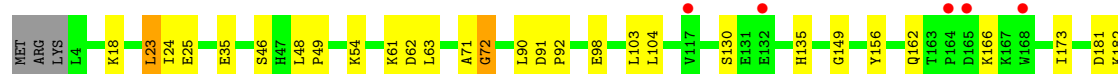
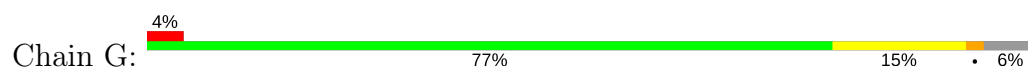
• Molecule 1: Ferulic acid decarboxylase 1

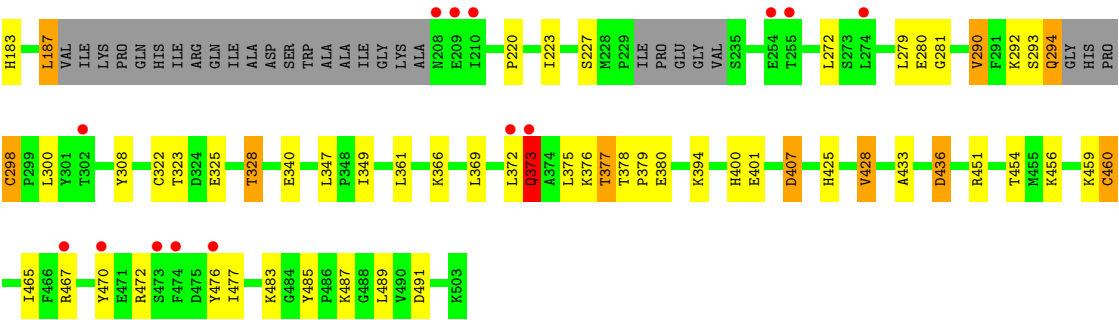


• Molecule 1: Ferulic acid decarboxylase 1

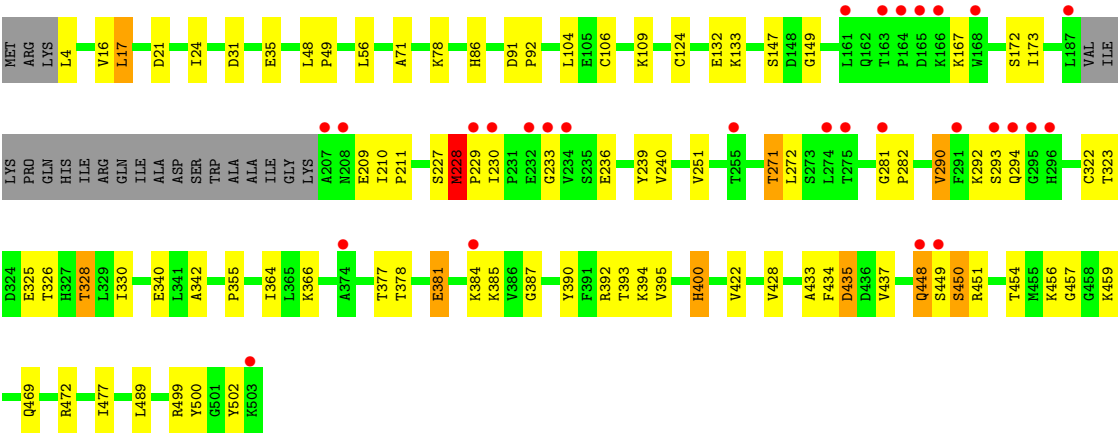
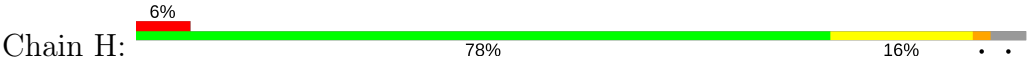


• Molecule 1: Ferulic acid decarboxylase 1





• Molecule 1: Ferulic acid decarboxylase 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	251.96Å 121.00Å 159.56Å 90.00° 121.93° 90.00°	Depositor
Resolution (Å)	47.05 – 2.35 49.85 – 2.35	Depositor EDS
% Data completeness (in resolution range)	95.8 (47.05-2.35) 91.5 (49.85-2.35)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.171 , 0.215 0.172 , 0.215	Depositor DCC
R_{free} test set	8147 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.559	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	31691	wwPDB-VP
Average B, all atoms (Å ²)	44.0	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 47.95 % 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 9.2275e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 4VP

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	0.43	0/3794	0.55	0/5147
1	B	0.43	0/3846	0.56	0/5218
1	C	0.41	0/4033	0.54	0/5477
1	D	0.45	0/3872	0.56	0/5256
1	E	0.41	0/3963	0.54	0/5376
1	F	0.41	0/4033	0.55	0/5477
1	G	0.39	0/3811	0.54	0/5170
1	H	0.39	0/3875	0.55	0/5262
All	All	0.42	0/31227	0.55	0/42383

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	C	0	1
1	G	0	3
1	H	0	1
All	All	0	5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	228	MET	Peptide
1	G	373	GLN	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	G	476	TYR	Peptide
1	G	72	GLY	Peptide
1	H	228	MET	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	3705	0	3702	35	0
1	B	3748	0	3761	45	0
1	C	3933	0	3944	54	1
1	D	3775	0	3781	35	0
1	E	3868	0	3882	40	0
1	F	3933	0	3944	57	0
1	G	3718	0	3722	68	1
1	H	3777	0	3776	69	0
2	A	9	0	8	2	0
2	C	9	0	8	2	0
2	E	9	0	8	1	0
2	F	9	0	8	0	0
2	G	9	0	8	0	0
3	A	152	0	0	8	0
3	B	190	0	0	9	0
3	C	131	0	0	7	0
3	D	185	0	0	11	0
3	E	165	0	0	10	0
3	F	137	0	0	6	0
3	G	100	0	0	7	0
3	H	129	0	0	9	0
All	All	31691	0	30552	382	1

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.

The worst 5 of 382 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:D:208:ASN:N	3:D:762:HOH:O	1.93	1.00
1:C:410:ASP:OD2	3:C:736:HOH:O	1.82	0.95
1:G:24[B]:ILE:HD11	1:G:49:PRO:HB2	1.55	0.88
1:C:228:MET:HB3	1:C:229:PRO:HD3	1.55	0.88
1:D:296:HIS:N	3:D:760:HOH:O	2.05	0.88

All (1) 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)
1:C:167:LYS:NZ	1:G:98:GLU:OE2[2_756]	2.15	0.05

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	464/503 (92%)	448 (97%)	15 (3%)	1 (0%)	51	59
1	B	470/503 (93%)	453 (96%)	17 (4%)	0	100	100
1	C	499/503 (99%)	476 (95%)	21 (4%)	2 (0%)	38	42
1	D	475/503 (94%)	455 (96%)	19 (4%)	1 (0%)	51	59
1	E	486/503 (97%)	472 (97%)	14 (3%)	0	100	100
1	F	499/503 (99%)	482 (97%)	16 (3%)	1 (0%)	51	59
1	G	465/503 (92%)	445 (96%)	18 (4%)	2 (0%)	38	42
1	H	478/503 (95%)	453 (95%)	23 (5%)	2 (0%)	38	42
All	All	3836/4024 (95%)	3684 (96%)	143 (4%)	9 (0%)	51	59

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	228	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	F	294	GLN
1	G	436	ASP
1	H	228	MET
1	C	293	SER

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	413/438 (94%)	404 (98%)	9 (2%)	57	70
1	B	420/438 (96%)	409 (97%)	11 (3%)	51	63
1	C	436/438 (100%)	419 (96%)	17 (4%)	37	48
1	D	422/438 (96%)	402 (95%)	20 (5%)	30	38
1	E	429/438 (98%)	421 (98%)	8 (2%)	62	74
1	F	436/438 (100%)	425 (98%)	11 (2%)	53	64
1	G	415/438 (95%)	398 (96%)	17 (4%)	35	45
1	H	421/438 (96%)	405 (96%)	16 (4%)	38	48
All	All	3392/3504 (97%)	3283 (97%)	109 (3%)	44	54

5 of 109 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	251	VAL
1	E	275	THR
1	H	290	VAL
1	D	271	THR
1	D	393	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	162	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	F	468	GLN
1	G	183	HIS
1	G	373	GLN
1	H	400	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](#)

5 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	4VP	A	601	-	9,9,9	0.63	0	11,11,11	0.92	0
2	4VP	C	601	-	9,9,9	0.64	0	11,11,11	0.86	0
2	4VP	E	601	-	9,9,9	0.77	0	11,11,11	0.76	0
2	4VP	F	601	-	9,9,9	0.77	0	11,11,11	0.67	0
2	4VP	G	601	-	9,9,9	0.72	0	11,11,11	0.77	0

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	4VP	A	601	-	-	0/2/2/2	0/1/1/1
2	4VP	C	601	-	-	0/2/2/2	0/1/1/1
2	4VP	E	601	-	-	0/2/2/2	0/1/1/1
2	4VP	F	601	-	-	0/2/2/2	0/1/1/1
2	4VP	G	601	-	-	0/2/2/2	0/1/1/1

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.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	4VP	2	0
2	C	601	4VP	2	0
2	E	601	4VP	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, 95th 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 ²)	Q<0.9
1	A	472/503 (93%)	-0.24	11 (2%) 61 70	22, 37, 66, 87	0
1	B	476/503 (94%)	-0.18	15 (3%) 48 59	23, 36, 66, 110	0
1	C	501/503 (99%)	-0.36	3 (0%) 89 93	25, 40, 64, 87	0
1	D	480/503 (95%)	-0.24	13 (2%) 55 64	19, 35, 66, 95	0
1	E	492/503 (97%)	-0.36	1 (0%) 94 97	24, 37, 62, 82	0
1	F	501/503 (99%)	-0.15	12 (2%) 59 68	22, 46, 75, 94	0
1	G	472/503 (93%)	0.05	19 (4%) 39 50	26, 53, 85, 110	0
1	H	481/503 (95%)	0.05	28 (5%) 24 34	26, 46, 84, 107	0
All	All	3875/4024 (96%)	-0.18	102 (2%) 56 65	19, 40, 74, 110	0

The worst 5 of 102 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	168	TRP	9.3
1	H	168	TRP	7.3
1	D	296	HIS	5.6
1	A	293	SER	5.1
1	B	167	LYS	5.0

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, 95th 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 ²)	Q<0.9
2	4VP	F	601	9/9	0.90	0.19	3.49	42,45,53,57	0
2	4VP	G	601	9/9	0.87	0.18	3.49	49,51,61,66	0
2	4VP	A	601	9/9	0.83	0.20	3.38	55,58,64,66	0
2	4VP	E	601	9/9	0.90	0.19	1.90	40,41,62,68	0
2	4VP	C	601	9/9	0.91	0.17	0.86	48,50,54,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.