



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 11:37 AM GMT

PDB ID : 3PCI
Title : STRUCTURE OF PROTOCATECHUATE 3,4-DIOXYGENASE COM-
PLEXED WITH 3-iodo-4-HYDROXYBENZOATE
Authors : Orville, A.M.; Elango, N.; Lipscomb, J.D.; Ohlendorf, D.H.
Deposited on : 1997-07-02
Resolution : 2.21 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 (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

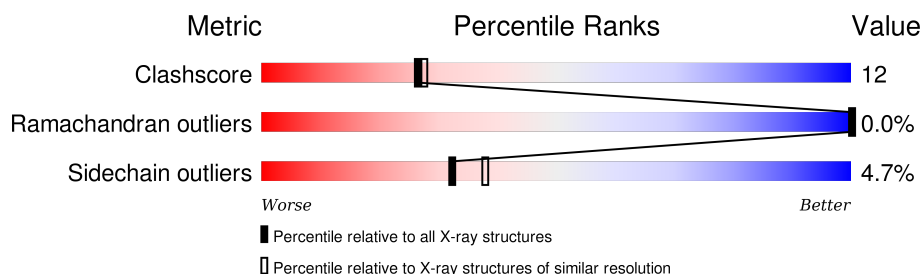
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	200	
1	B	200	
1	C	200	
1	D	200	
1	E	200	
1	F	200	
2	M	238	

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Mol	Chain	Length	Quality of chain
2	N	238	<div><div></div><div>76%</div><div>18%</div><div></div><div>• • •</div></div>
2	O	238	<div><div></div><div>76%</div><div>19%</div><div></div><div>• •</div></div>
2	P	238	<div><div></div><div>77%</div><div>17%</div><div></div><div>• •</div></div>
2	Q	238	<div><div></div><div>69%</div><div>21%</div><div>7%</div><div>•</div></div>
2	R	238	<div><div></div><div>68%</div><div>24%</div><div>5%</div><div>• •</div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 22104 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 PROTOCATECHUATE 3,4-DIOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	B	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	C	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	D	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	E	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			
1	F	200	Total	C	N	O	S	0	0	0
			1571	993	276	299	3			

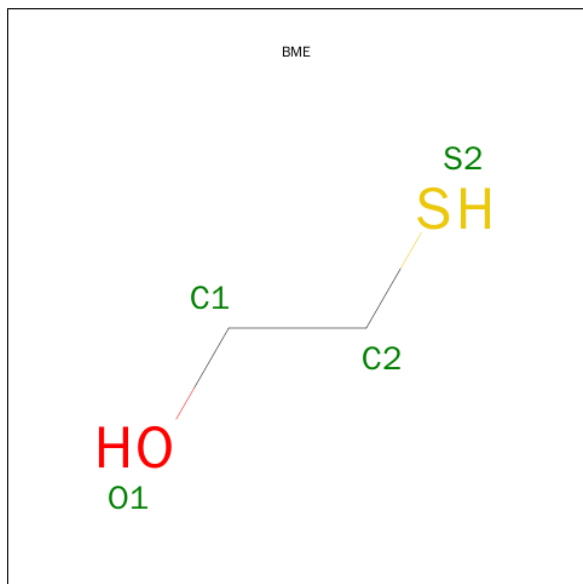
- Molecule 2 is a protein called PROTOCATECHUATE 3,4-DIOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	233	Total	C	N	O	S	0	0	0
			1840	1171	334	328	7			
2	N	233	Total	C	N	O	S	0	0	0
			1840	1171	334	328	7			
2	O	233	Total	C	N	O	S	0	0	0
			1840	1171	334	328	7			
2	P	233	Total	C	N	O	S	0	0	0
			1840	1171	334	328	7			
2	Q	233	Total	C	N	O	S	0	0	0
			1840	1171	334	328	7			
2	R	233	Total	C	N	O	S	0	0	0
			1840	1171	334	328	7			

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe).

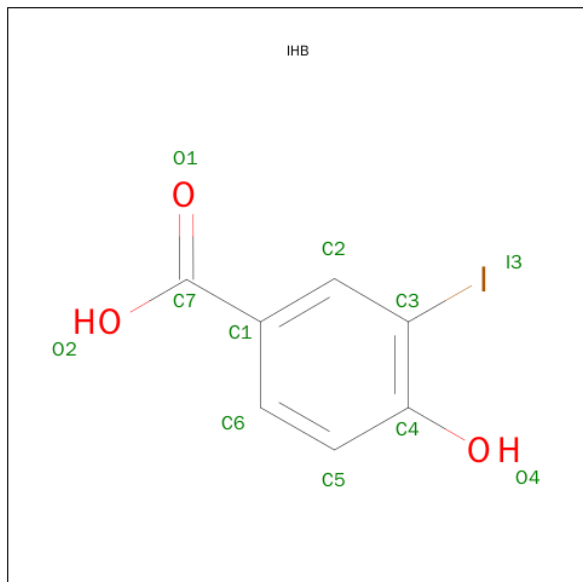
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	P	1	Total Fe 1 1	0	0
3	Q	1	Total Fe 1 1	0	0
3	N	1	Total Fe 1 1	0	0
3	O	1	Total Fe 1 1	0	0
3	R	1	Total Fe 1 1	0	0
3	M	1	Total Fe 1 1	0	0

- Molecule 4 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	M	1	Total C O S 4 2 1 1	0	0
4	N	1	Total C O S 4 2 1 1	0	0
4	O	1	Total C O S 4 2 1 1	0	0
4	P	1	Total C O S 4 2 1 1	0	0
4	Q	1	Total C O S 4 2 1 1	0	0
4	R	1	Total C O S 4 2 1 1	0	0

- Molecule 5 is 3-iodo-4-hydroxybenzoic acid (three-letter code: IHB) (formula: $C_7H_5IO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	M	1	Total	C	I	O	0	0
			11	7	1	3		
5	M	1	Total	C	I	O	0	0
			11	7	1	3		
5	N	1	Total	C	I	O	0	0
			11	7	1	3		
5	N	1	Total	C	I	O	0	0
			11	7	1	3		
5	O	1	Total	C	I	O	0	0
			11	7	1	3		
5	O	1	Total	C	I	O	0	0
			11	7	1	3		
5	P	1	Total	C	I	O	0	0
			11	7	1	3		
5	P	1	Total	C	I	O	0	0
			11	7	1	3		
5	Q	1	Total	C	I	O	0	0
			11	7	1	3		
5	Q	1	Total	C	I	O	0	0
			11	7	1	3		
5	R	1	Total	C	I	O	0	0
			11	7	1	3		
5	R	1	Total	C	I	O	0	0
			11	7	1	3		

- Molecule 6 is water.

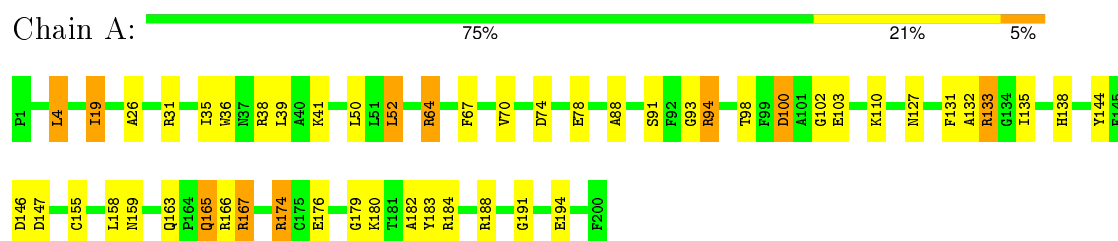
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	82	Total 82	O 82	0	0
6	B	85	Total 85	O 85	0	0
6	C	82	Total 82	O 82	0	0
6	D	81	Total 81	O 81	0	0
6	E	84	Total 84	O 84	0	0
6	F	81	Total 81	O 81	0	0
6	M	163	Total 163	O 163	0	0
6	N	167	Total 167	O 167	0	0
6	O	162	Total 162	O 162	0	0
6	P	157	Total 157	O 157	0	0
6	Q	165	Total 165	O 165	0	0
6	R	167	Total 167	O 167	0	0

3 Residue-property plots

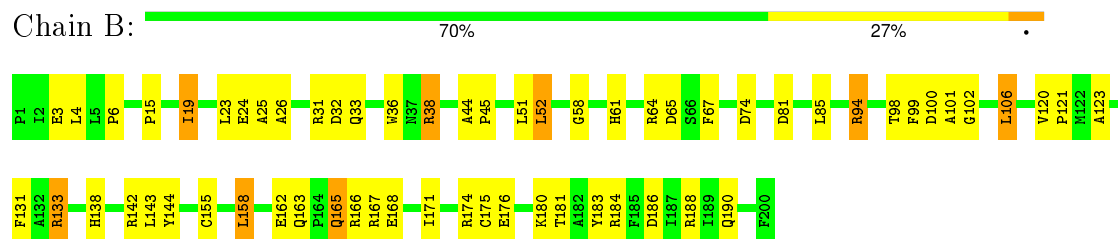
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

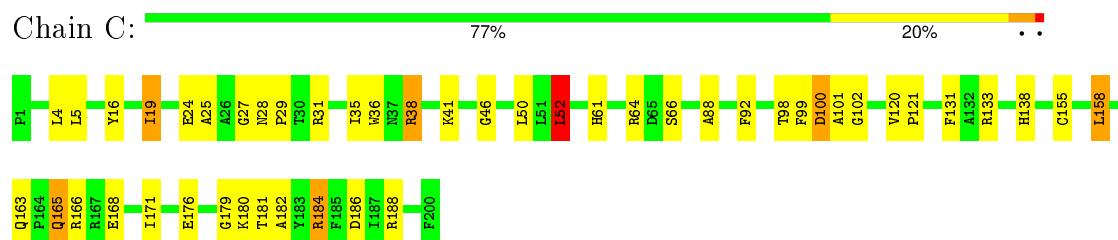
• Molecule 1: PROTOCATECHUATE 3,4-DIOXYGENASE



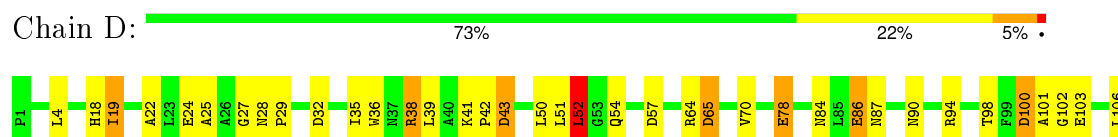
• Molecule 1: PROTOCATECHUATE 3,4-DIOXYGENASE

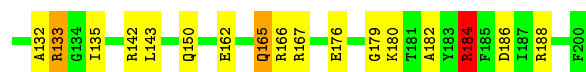


• Molecule 1: PROTOCATECHUATE 3,4-DIOXYGENASE



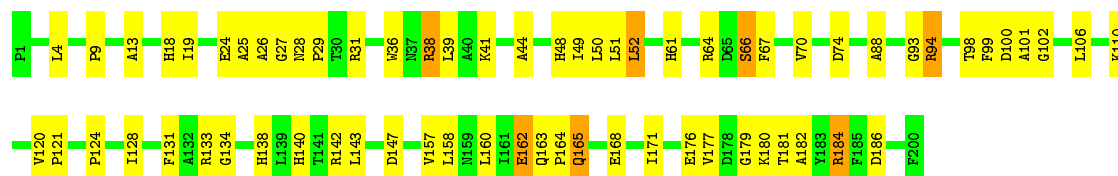
• Molecule 1: PROTOCATECHUATE 3,4-DIOXYGENASE





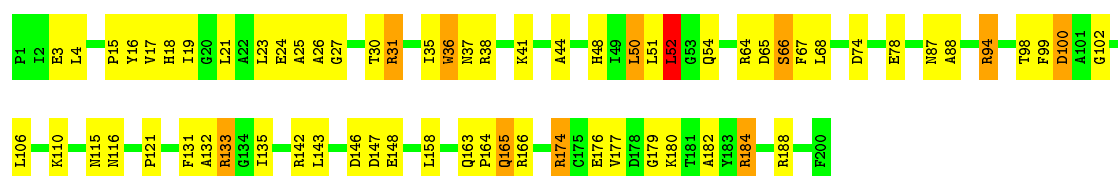
• Molecule 1: PROTOCATECHUATE 3,4-DIOXYGENASE

Chain E: 67% 30%



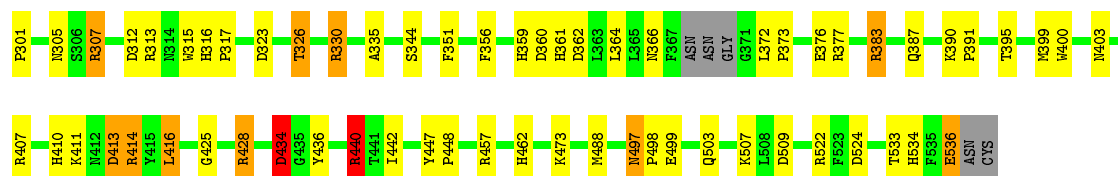
• Molecule 1: PROTOCATECHUATE 3,4-DIOXYGENASE

Chain F: 67% 28% 5%



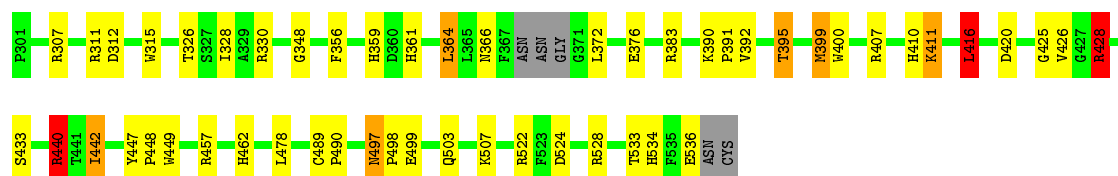
• Molecule 2: PROTOCATECHUATE 3,4-DIOXYGENASE

Chain M: 72% 21%



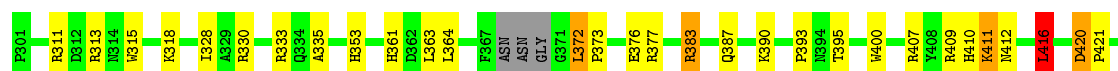
• Molecule 2: PROTOCATECHUATE 3,4-DIOXYGENASE

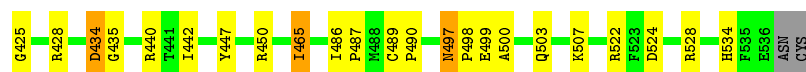
Chain N: 76% 18%



• Molecule 2: PROTOCATECHUATE 3,4-DIOXYGENASE

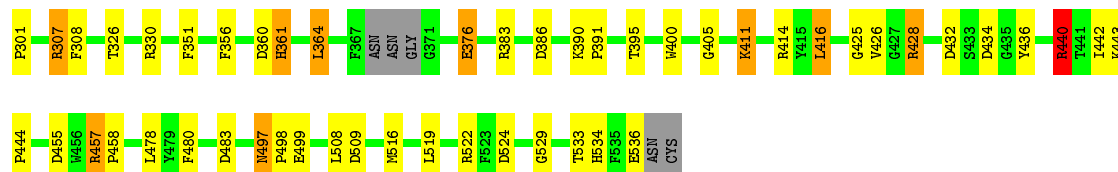
Chain O: 76% 19%





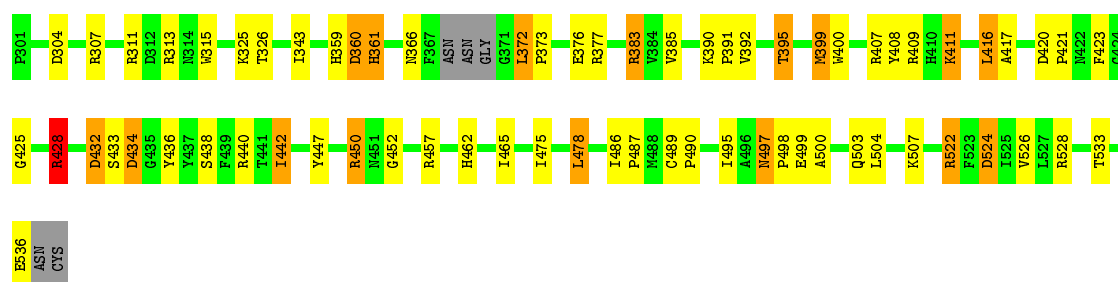
• Molecule 2: PROTOCATECHUATE 3,4-DIOXYGENASE

Chain P: 77% 17%



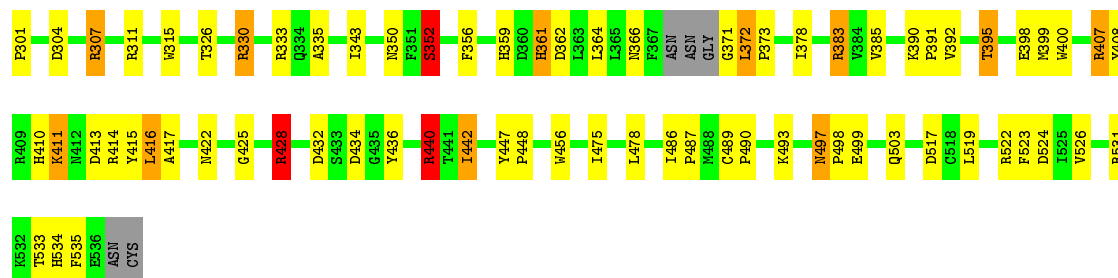
• Molecule 2: PROTOCATECHUATE 3,4-DIOXYGENASE

Chain Q: 69% 21% 7%



• Molecule 2: PROTOCATECHUATE 3,4-DIOXYGENASE

Chain R: 68% 24% 5%



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	197.00Å 127.07Å 134.08Å 90.00° 97.53° 90.00°	Depositor
Resolution (Å)	6.00 – 2.21	Depositor
% Data completeness (in resolution range)	83.0 (6.00-2.21)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.159 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	22104	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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: IHB, FE, BME

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	0.94	1/1611 (0.1%)	1.66	30/2195 (1.4%)
1	B	0.95	0/1611	1.66	33/2195 (1.5%)
1	C	0.95	1/1611 (0.1%)	1.50	14/2195 (0.6%)
1	D	0.92	0/1611	1.52	23/2195 (1.0%)
1	E	0.97	0/1611	1.52	23/2195 (1.0%)
1	F	0.98	0/1611	1.55	25/2195 (1.1%)
2	M	1.00	0/1895	1.54	25/2580 (1.0%)
2	N	0.98	0/1895	1.50	16/2580 (0.6%)
2	O	1.00	0/1895	1.55	24/2580 (0.9%)
2	P	0.97	0/1895	1.53	21/2580 (0.8%)
2	Q	1.02	1/1895 (0.1%)	1.54	24/2580 (0.9%)
2	R	1.00	1/1895 (0.1%)	1.56	24/2580 (0.9%)
All	All	0.98	4/21036 (0.0%)	1.55	282/28650 (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	D	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	ARG	CD-NE	-5.16	1.37	1.46
1	C	66	SER	CB-OG	-5.10	1.35	1.42
2	R	428	ARG	CD-NE	-5.07	1.37	1.46
2	Q	433	SER	CB-OG	5.05	1.48	1.42

The worst 5 of 282 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	440	ARG	NE-CZ-NH2	-20.04	110.28	120.30
2	P	440	ARG	NE-CZ-NH2	-19.34	110.63	120.30
2	R	440	ARG	NE-CZ-NH2	-18.26	111.17	120.30
2	M	440	ARG	NE-CZ-NH2	-16.43	112.08	120.30
2	Q	440	ARG	NE-CZ-NH2	-15.94	112.33	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	184	ARG	Sidechain

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	1571	0	1499	37	0
1	B	1571	0	1499	40	0
1	C	1571	0	1499	29	1
1	D	1571	0	1499	35	1
1	E	1571	0	1499	47	0
1	F	1571	0	1499	66	0
2	M	1840	0	1792	57	0
2	N	1840	0	1792	39	0
2	O	1840	0	1792	36	0
2	P	1840	0	1792	41	0
2	Q	1840	0	1792	56	0
2	R	1840	0	1792	56	0
3	M	1	0	0	0	0
3	N	1	0	0	0	0
3	O	1	0	0	0	0
3	P	1	0	0	0	0
3	Q	1	0	0	0	0
3	R	1	0	0	0	0
4	M	4	0	5	1	0
4	N	4	0	5	0	0
4	O	4	0	5	1	0
4	P	4	0	5	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	Q	4	0	5	2	0
4	R	4	0	5	0	0
5	M	22	0	7	1	0
5	N	22	0	7	1	0
5	O	22	0	7	2	0
5	P	22	0	7	0	0
5	Q	22	0	7	2	0
5	R	22	0	7	3	0
6	A	82	0	0	0	0
6	B	85	0	0	0	0
6	C	82	0	0	0	0
6	D	81	0	0	0	0
6	E	84	0	0	2	0
6	F	81	0	0	2	0
6	M	163	0	0	7	0
6	N	167	0	0	3	0
6	O	162	0	0	6	0
6	P	157	0	0	4	0
6	Q	165	0	0	8	1
6	R	167	0	0	4	0
All	All	22104	0	19818	492	2

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 12.

The worst 5 of 492 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)
2:M:364:LEU:HD22	2:M:440:ARG:HD3	1.25	1.16
2:P:364:LEU:HD22	2:P:440:ARG:HD3	1.22	1.12
1:D:64:ARG:NH1	1:D:100:ASP:O	1.81	1.12
1:B:165:GLN:NE2	1:B:165:GLN:H	1.49	1.10
1:F:165:GLN:NE2	1:F:165:GLN:H	1.50	1.09

All (2) 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:46:GLY:O	1:D:150:GLN:OE1[3_554]	1.84	0.36
6:Q:1200:HOH:O	6:Q:1200:HOH:O[2_555]	2.10	0.10

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	198/200 (99%)	192 (97%)	6 (3%)	0	100	100
1	B	198/200 (99%)	192 (97%)	6 (3%)	0	100	100
1	C	198/200 (99%)	193 (98%)	5 (2%)	0	100	100
1	D	198/200 (99%)	189 (96%)	9 (4%)	0	100	100
1	E	198/200 (99%)	193 (98%)	5 (2%)	0	100	100
1	F	198/200 (99%)	192 (97%)	6 (3%)	0	100	100
2	M	229/238 (96%)	222 (97%)	7 (3%)	0	100	100
2	N	229/238 (96%)	220 (96%)	9 (4%)	0	100	100
2	O	229/238 (96%)	221 (96%)	8 (4%)	0	100	100
2	P	229/238 (96%)	222 (97%)	7 (3%)	0	100	100
2	Q	229/238 (96%)	218 (95%)	11 (5%)	0	100	100
2	R	229/238 (96%)	221 (96%)	7 (3%)	1 (0%)	39	41
All	All	2562/2628 (98%)	2475 (97%)	86 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	R	535	PHE

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	162/163 (99%)	156 (96%)	6 (4%)	41	50
1	B	162/163 (99%)	155 (96%)	7 (4%)	35	42
1	C	162/163 (99%)	155 (96%)	7 (4%)	35	42
1	D	162/163 (99%)	155 (96%)	7 (4%)	35	42
1	E	162/163 (99%)	155 (96%)	7 (4%)	35	42
1	F	162/163 (99%)	155 (96%)	7 (4%)	35	42
2	M	196/202 (97%)	188 (96%)	8 (4%)	37	45
2	N	196/202 (97%)	183 (93%)	13 (7%)	21	21
2	O	196/202 (97%)	186 (95%)	10 (5%)	29	33
2	P	196/202 (97%)	189 (96%)	7 (4%)	42	51
2	Q	196/202 (97%)	185 (94%)	11 (6%)	26	29
2	R	196/202 (97%)	185 (94%)	11 (6%)	26	29
All	All	2148/2190 (98%)	2047 (95%)	101 (5%)	32	38

5 of 101 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	O	416	LEU
1	D	52	LEU
2	R	399	MET
2	O	465	ILE
1	D	4	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 34 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	163	GLN
2	P	497	ASN
2	R	497	ASN
2	P	361	HIS
1	B	165	GLN

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 24 ligands modelled in this entry, 6 are monoatomic - leaving 18 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	$\# Z > 2$
5	IHB	M	550	3	8,11,11	1.33	1 (12%)	10,15,15	0.65	0
5	IHB	M	551	-	8,11,11	1.20	1 (12%)	10,15,15	0.55	0
4	BME	M	601	2	3,3,3	0.55	0	2,2,2	0.12	0
5	IHB	N	550	3	8,11,11	1.22	1 (12%)	10,15,15	0.99	0
5	IHB	N	551	-	8,11,11	0.89	0	10,15,15	0.68	0
4	BME	N	601	2	3,3,3	0.50	0	2,2,2	0.07	0
5	IHB	O	550	3	8,11,11	1.00	1 (12%)	10,15,15	0.53	0
5	IHB	O	551	-	8,11,11	1.28	1 (12%)	10,15,15	0.89	0
4	BME	O	601	2	3,3,3	0.66	0	2,2,2	0.29	0
5	IHB	P	550	3	8,11,11	1.18	1 (12%)	10,15,15	0.57	0
5	IHB	P	551	-	8,11,11	1.21	1 (12%)	10,15,15	0.82	0
4	BME	P	601	2	3,3,3	0.54	0	2,2,2	0.56	0
5	IHB	Q	550	3	8,11,11	1.16	1 (12%)	10,15,15	0.53	0
5	IHB	Q	551	-	8,11,11	0.87	1 (12%)	10,15,15	0.65	0
4	BME	Q	601	2	3,3,3	0.35	0	2,2,2	0.55	0
5	IHB	R	550	3	8,11,11	1.28	1 (12%)	10,15,15	0.55	0
5	IHB	R	551	-	8,11,11	1.13	1 (12%)	10,15,15	0.84	0
4	BME	R	601	2	3,3,3	0.52	0	2,2,2	0.41	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
5	IHB	M	550	3	-	0/0/4/4	0/1/1/1
5	IHB	M	551	-	-	0/0/4/4	0/1/1/1
4	BME	M	601	2	-	0/1/1/1	0/0/0/0
5	IHB	N	550	3	-	0/0/4/4	0/1/1/1
5	IHB	N	551	-	-	0/0/4/4	0/1/1/1
4	BME	N	601	2	-	0/1/1/1	0/0/0/0
5	IHB	O	550	3	-	0/0/4/4	0/1/1/1
5	IHB	O	551	-	-	0/0/4/4	0/1/1/1
4	BME	O	601	2	-	0/1/1/1	0/0/0/0
5	IHB	P	550	3	-	0/0/4/4	0/1/1/1
5	IHB	P	551	-	-	0/0/4/4	0/1/1/1
4	BME	P	601	2	-	0/1/1/1	0/0/0/0
5	IHB	Q	550	3	-	0/0/4/4	0/1/1/1
5	IHB	Q	551	-	-	0/0/4/4	0/1/1/1
4	BME	Q	601	2	-	0/1/1/1	0/0/0/0
5	IHB	R	550	3	-	0/0/4/4	0/1/1/1
5	IHB	R	551	-	-	0/0/4/4	0/1/1/1
4	BME	R	601	2	-	0/1/1/1	0/0/0/0

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	M	550	IHB	C3-I3	-3.43	2.01	2.10
5	R	550	IHB	C3-I3	-3.23	2.01	2.10
5	O	551	IHB	C3-I3	-3.18	2.01	2.10
5	N	550	IHB	C3-I3	-3.01	2.02	2.10
5	P	551	IHB	C3-I3	-2.92	2.02	2.10

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	M	550	IHB	1	0
4	M	601	BME	1	0
5	N	550	IHB	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	O	550	IHB	2	0
4	O	601	BME	1	0
5	Q	550	IHB	2	0
4	Q	601	BME	2	0
5	R	550	IHB	3	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](#)

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

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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