



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 08:32 PM UTC

PDB ID : 3PCF / pdb_00003pcf
Title : STRUCTURE OF PROTOCATECHUATE 3,4-DIOXYGENASE COM-
PLEXED WITH 3-FLURO-4-HYDROXYBENZOATE
Authors : Orville, A.M.; Elango, N.; Lipscomb, J.D.; Ohlendorf, D.H.
Deposited on : 1997-06-27
Resolution : 2.15 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

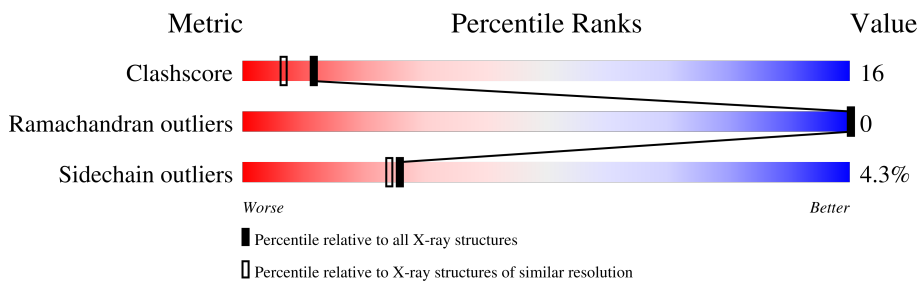
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)





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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$

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	
2	N	238	

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Mol	Chain	Length	Quality of chain
2	O	238	 63% 29% 5% •
2	P	238	 61% 31% 6% •
2	Q	238	 59% 32% 7% •
2	R	238	 58% 32% 6% ••

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 22080 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 alpha chain.

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

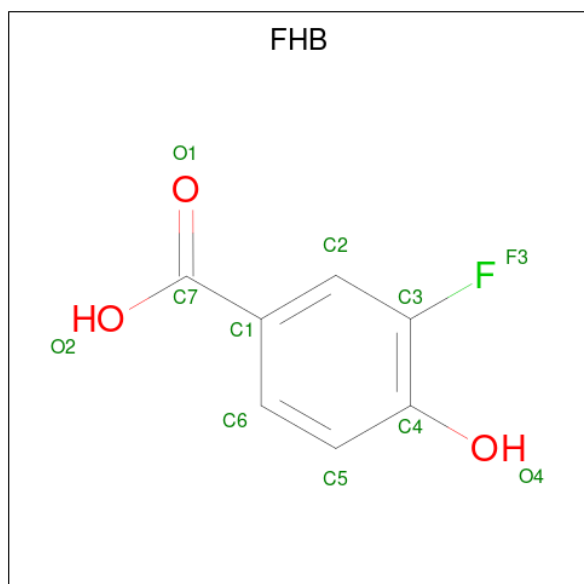
- Molecule 2 is a protein called Protocatechuate 3,4-dioxygenase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	M	233	1844	1173	334	329	8	0	0	0
2	N	233	1844	1173	334	329	8	0	0	0
2	O	233	1844	1173	334	329	8	0	0	0
2	P	233	1844	1173	334	329	8	0	0	0
2	Q	233	1844	1173	334	329	8	0	0	0
2	R	233	1844	1173	334	329	8	0	0	0

- Molecule 3 is FE (III) ION (CCD ID: FE) (formula: Fe).

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

- Molecule 4 is 3-FLUORO-4-HYDROXYBENZOIC ACID (CCD ID: FHB) (formula: $C_7H_5FO_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	M	1	Total C F O 11 7 1 3	0	0
4	M	1	Total C F O 11 7 1 3	0	0
4	N	1	Total C F O 11 7 1 3	0	0
4	N	1	Total C F O 11 7 1 3	0	0
4	O	1	Total C F O 11 7 1 3	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	O	1	Total	C	F	O	0	0
			11	7	1	3		
4	P	1	Total	C	F	O	0	0
			11	7	1	3		
4	P	1	Total	C	F	O	0	0
			11	7	1	3		
4	Q	1	Total	C	F	O	0	0
			11	7	1	3		
4	Q	1	Total	C	F	O	0	0
			11	7	1	3		
4	R	1	Total	C	F	O	0	0
			11	7	1	3		
4	R	1	Total	C	F	O	0	0
			11	7	1	3		

- Molecule 5 is water.

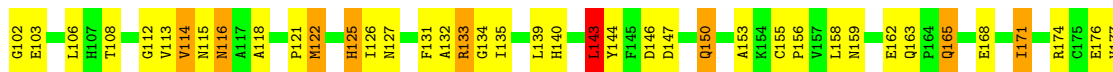
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	85	Total	O	0	0
			85	85		
5	M	158	Total	O	0	0
			158	158		
5	B	80	Total	O	0	0
			80	80		
5	N	166	Total	O	0	0
			166	166		
5	C	86	Total	O	0	0
			86	86		
5	O	155	Total	O	0	0
			155	155		
5	D	80	Total	O	0	0
			80	80		
5	P	155	Total	O	0	0
			155	155		
5	E	84	Total	O	0	0
			84	84		
5	Q	159	Total	O	0	0
			159	159		
5	F	83	Total	O	0	0
			83	83		
5	R	161	Total	O	0	0
			161	161		



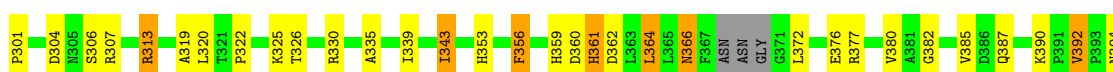
- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain



- Molecule 1: Protocatechuate 3,4-dioxygenase alpha chain

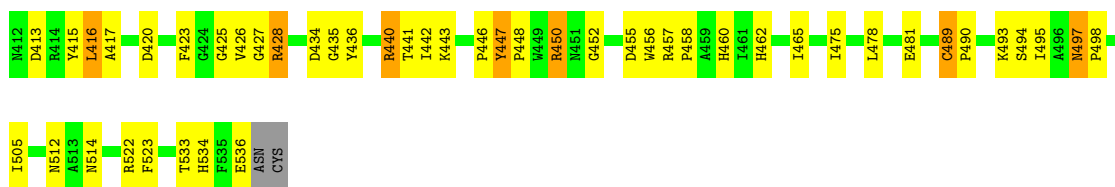


- Molecule 2: Protocatechuate 3,4-dioxygenase beta chain



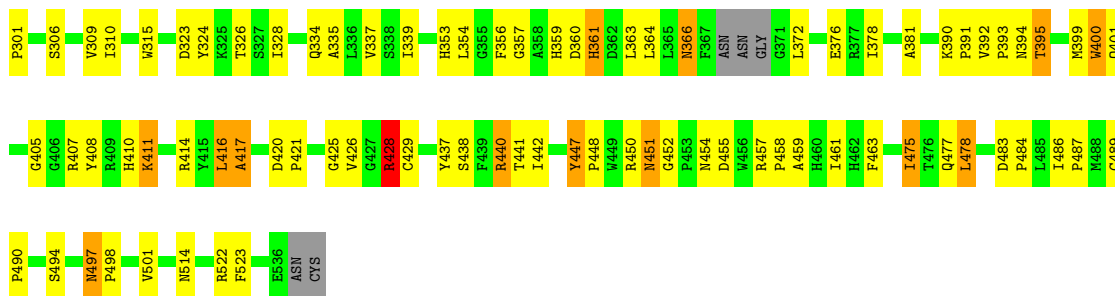
- Molecule 2: Protocatechuate 3,4-dioxygenase beta chain





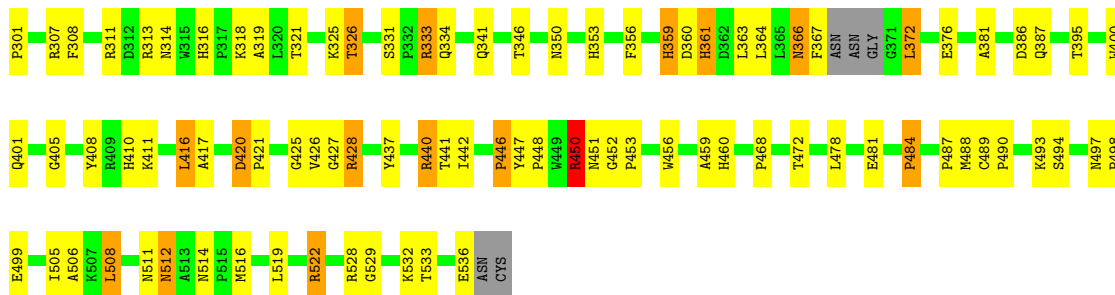
- Molecule 2: Protocatechuic 3,4-dioxygenase beta chain

Chain O: 63% 29% 5%



- Molecule 2: Protocatechuic 3,4-dioxygenase beta chain

Chain P: 61% 31% 6%



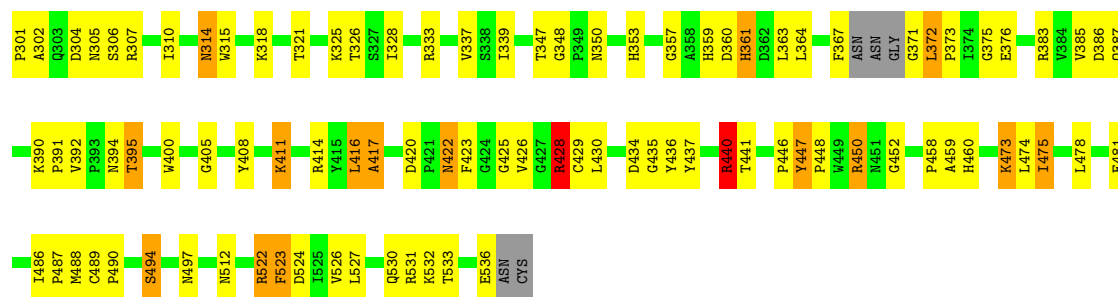
- Molecule 2: Protocatechuic 3,4-dioxygenase beta chain

Chain Q: 59% 32% 7%



- Molecule 2: Protocatechuic 3,4-dioxygenase beta chain

Chain R: 58% 32% 6%



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	196.03Å 127.22Å 133.70Å 90.00° 97.70° 90.00°	Depositor
Resolution (Å)	6.00 – 2.15	Depositor
% Data completeness (in resolution range)	93.3 (6.00-2.15)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.179 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	22080	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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: FE, CME, FHB

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	1.29	2/1611 (0.1%)	1.93	36/2195 (1.6%)
1	B	1.29	4/1611 (0.2%)	1.82	26/2195 (1.2%)
1	C	1.32	3/1611 (0.2%)	1.81	29/2195 (1.3%)
1	D	1.33	2/1611 (0.1%)	1.88	32/2195 (1.5%)
1	E	1.37	2/1611 (0.1%)	1.84	30/2195 (1.4%)
1	F	1.37	4/1611 (0.2%)	1.86	38/2195 (1.7%)
2	M	1.48	9/1888 (0.5%)	2.00	53/2569 (2.1%)
2	N	1.40	4/1888 (0.2%)	1.90	32/2569 (1.2%)
2	O	1.45	7/1888 (0.4%)	1.93	41/2569 (1.6%)
2	P	1.47	8/1888 (0.4%)	1.95	54/2569 (2.1%)
2	Q	1.42	5/1888 (0.3%)	1.93	48/2569 (1.9%)
2	R	1.43	3/1888 (0.2%)	1.93	42/2569 (1.6%)
All	All	1.39	53/20994 (0.3%)	1.90	461/28584 (1.6%)

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

The worst 5 of 53 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	ARG	CD-NE	-9.96	1.32	1.46
2	P	451	ASN	CA-C	9.88	1.57	1.52
2	O	451	ASN	CA-C	9.46	1.56	1.52
2	M	451	ASN	CA-C	8.71	1.56	1.52
2	P	428	ARG	CD-NE	-8.07	1.34	1.46

The worst 5 of 461 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	ARG	CD-NE-CZ	22.86	156.41	124.40
1	D	133	ARG	CD-NE-CZ	15.40	145.96	124.40
2	Q	450	ARG	CD-NE-CZ	13.79	143.71	124.40
2	P	428	ARG	CG-CD-NE	12.06	138.53	112.00
2	P	440	ARG	NE-CZ-NH2	-10.89	109.40	119.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	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	47	0
1	B	1571	0	1499	51	0
1	C	1571	0	1499	66	0
1	D	1571	0	1499	59	0
1	E	1571	0	1499	62	0
1	F	1571	0	1499	82	0
2	M	1844	0	1796	62	0
2	N	1844	0	1796	54	0
2	O	1844	0	1796	46	0
2	P	1844	0	1796	61	0
2	Q	1844	0	1796	67	0
2	R	1844	0	1796	71	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	22	0	7	0	0
4	N	22	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	O	22	0	7	2	0
4	P	22	0	7	0	0
4	Q	22	0	7	2	0
4	R	22	0	7	2	0
5	A	85	0	0	3	0
5	B	80	0	0	2	0
5	C	86	0	0	1	0
5	D	80	0	0	1	0
5	E	84	0	0	3	0
5	F	83	0	0	2	0
5	M	158	0	0	5	0
5	N	166	0	0	4	0
5	O	155	0	0	4	0
5	P	155	0	0	3	0
5	Q	159	0	0	7	0
5	R	161	0	0	4	0
All	All	22080	0	19812	644	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 16.

The worst 5 of 644 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:64:ARG:NH1	1:D:100:ASP:O	1.87	1.07
1:F:64:ARG:NH1	1:F:100:ASP:O	1.87	1.06
1:B:165:GLN:H	1:B:165:GLN:NE2	1.58	1.00
1:E:165:GLN:H	1:E:165:GLN:NE2	1.59	0.97
1:C:64:ARG:NH1	1:C:100:ASP:O	1.97	0.96

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	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%)	188 (95%)	10 (5%)	0	100	100
1	F	198/200 (99%)	185 (93%)	13 (7%)	0	100	100
2	M	228/238 (96%)	218 (96%)	10 (4%)	0	100	100
2	N	228/238 (96%)	221 (97%)	7 (3%)	0	100	100
2	O	228/238 (96%)	220 (96%)	8 (4%)	0	100	100
2	P	228/238 (96%)	219 (96%)	9 (4%)	0	100	100
2	Q	228/238 (96%)	219 (96%)	9 (4%)	0	100	100
2	R	228/238 (96%)	218 (96%)	10 (4%)	0	100	100
All	All	2556/2628 (97%)	2454 (96%)	102 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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%)	155 (96%)	7 (4%)	26	24
1	B	162/163 (99%)	152 (94%)	10 (6%)	16	12
1	C	162/163 (99%)	154 (95%)	8 (5%)	22	19
1	D	162/163 (99%)	154 (95%)	8 (5%)	22	19
1	E	162/163 (99%)	153 (94%)	9 (6%)	19	15
1	F	162/163 (99%)	156 (96%)	6 (4%)	30	30
2	M	195/201 (97%)	185 (95%)	10 (5%)	21	18

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	N	195/201 (97%)	187 (96%)	8 (4%)	27	26
2	O	195/201 (97%)	186 (95%)	9 (5%)	24	22
2	P	195/201 (97%)	190 (97%)	5 (3%)	40	43
2	Q	195/201 (97%)	189 (97%)	6 (3%)	35	37
2	R	195/201 (97%)	188 (96%)	7 (4%)	31	31
All	All	2142/2184 (98%)	2049 (96%)	93 (4%)	26	24

5 of 93 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	126	ILE
1	E	165	GLN
1	D	165	GLN
1	E	4	LEU
2	Q	390	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 42 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	11	GLN
1	F	150	GLN
1	E	163	GLN
2	Q	530	GLN
2	R	361	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](#)

6 non-standard protein/DNA/RNA residues 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	CME	Q	429	2	8,9,10	0.81	0	6,9,11	1.25	1 (16%)
2	CME	P	429	2	8,9,10	0.84	0	6,9,11	1.03	0
2	CME	O	429	2	8,9,10	0.81	0	6,9,11	1.13	0
2	CME	M	429	2	8,9,10	0.80	0	6,9,11	0.66	0
2	CME	R	429	2	8,9,10	0.94	0	6,9,11	1.34	0
2	CME	N	429	2	8,9,10	0.74	0	6,9,11	0.61	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	CME	Q	429	2	-	1/5/8/10	-
2	CME	P	429	2	-	1/5/8/10	-
2	CME	O	429	2	-	1/5/8/10	-
2	CME	M	429	2	-	1/5/8/10	-
2	CME	R	429	2	-	1/5/8/10	-
2	CME	N	429	2	-	0/5/8/10	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	Q	429	CME	CZ-CE-SD	-2.36	105.51	113.39

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	R	429	CME	CZ-CE-SD-SG
2	M	429	CME	SD-CE-CZ-OH
2	O	429	CME	CZ-CE-SD-SG
2	Q	429	CME	CZ-CE-SD-SG
2	P	429	CME	CZ-CE-SD-SG

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	Q	429	CME	3	0
2	O	429	CME	2	0
2	R	429	CME	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 6 are monoatomic - leaving 12 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$
4	FHB	N	550	3	11,11,11	1.29	1 (9%)	13,15,15	0.75	0
4	FHB	M	550	3	11,11,11	1.33	1 (9%)	13,15,15	1.11	2 (15%)
4	FHB	Q	551	-	11,11,11	1.11	0	13,15,15	1.06	1 (7%)
4	FHB	P	551	-	11,11,11	1.26	1 (9%)	13,15,15	1.21	1 (7%)
4	FHB	O	551	-	11,11,11	1.13	0	13,15,15	0.93	0
4	FHB	M	551	-	11,11,11	1.26	1 (9%)	13,15,15	1.07	2 (15%)
4	FHB	Q	550	3	11,11,11	1.49	2 (18%)	13,15,15	0.98	0
4	FHB	N	551	-	11,11,11	1.14	1 (9%)	13,15,15	1.14	2 (15%)
4	FHB	R	551	-	11,11,11	1.10	0	13,15,15	1.01	1 (7%)
4	FHB	P	550	3	11,11,11	1.26	1 (9%)	13,15,15	1.14	2 (15%)
4	FHB	O	550	3	11,11,11	1.27	2 (18%)	13,15,15	0.94	0
4	FHB	R	550	3	11,11,11	1.06	1 (9%)	13,15,15	1.00	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
4	FHB	N	550	3	-	0/4/4/4	0/1/1/1
4	FHB	M	550	3	-	0/4/4/4	0/1/1/1
4	FHB	Q	551	-	-	0/4/4/4	0/1/1/1
4	FHB	P	551	-	-	0/4/4/4	0/1/1/1
4	FHB	O	551	-	-	0/4/4/4	0/1/1/1
4	FHB	M	551	-	-	0/4/4/4	0/1/1/1
4	FHB	Q	550	3	-	0/4/4/4	0/1/1/1
4	FHB	N	551	-	-	0/4/4/4	0/1/1/1
4	FHB	R	551	-	-	0/4/4/4	0/1/1/1
4	FHB	P	550	3	-	0/4/4/4	0/1/1/1
4	FHB	O	550	3	-	0/4/4/4	0/1/1/1
4	FHB	R	550	3	-	0/4/4/4	0/1/1/1

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	Q	550	FHB	C2-C1	2.93	1.43	1.39
4	M	550	FHB	O2-C7	-2.90	1.21	1.30
4	N	550	FHB	O2-C7	-2.80	1.22	1.30
4	O	550	FHB	C4-C3	2.71	1.41	1.39
4	Q	550	FHB	O2-C7	-2.67	1.22	1.30

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	551	FHB	O2-C7-C1	2.74	121.88	114.84
4	N	551	FHB	O2-C7-C1	2.56	121.41	114.84
4	P	550	FHB	O2-C7-C1	2.50	121.25	114.84
4	M	551	FHB	O2-C7-C1	2.45	121.12	114.84
4	R	550	FHB	O2-C7-O1	-2.34	118.33	123.35

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	Q	550	FHB	2	0
4	O	550	FHB	2	0
4	R	550	FHB	2	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.