The PubChem System generates a binary substructure fingerprint for chemical structures. These fingerprints are used by PubChem for similarity neighboring and similarity searching.

A substructure is a fragment of a chemical structure. A fingerprint is an ordered list of binary (1/0) bits. Each bit represents a Boolean determination of, or test for, the presence of, for example, an element count, a type of ring system, atom pairing, atom environment (nearest neighbors), etc., in a chemical structure.

The native format of the PubChem Substructure Fingerprint property is binary data with a four byte integer prefix, where this integer prefix indicates the length of the bit list. For the ASN.1 and XML formatted data, this property is stored in a PC-InfoData container, as described by the PCSubstance ASN.1 definition or XML schema: ftp://ftp.ncbi.nlm.nih.gov/pubchem/specifications/

PC-InfoData is able to handle various types of data. Each PC-InfoData has a PC-Urn object (urn = universal resource name). Each property has a unique trio of "label", "name", and "datatype" definition (e.g., for PubChem Substructure Fingerprint, this is "Fingerprint", "SubStructure Keys", and "fingerprint", respectively). The fingerprint binary data is hex-encoded, when provided in the XML or textual ASN.1 formats.

When exporting fingerprint information in the SD file format, the SD tag for the PubChem Substructure Fingerprint property is "PUBCHEM\_CACTVS\_SUBGRAPHKEYS". The PubChem Substructure Fingerprint is Base64 encoded to provide a textual representation of the binary data. For a description of the Base64 encoding and decoding algorithm specification, go to: http://www.fags.org/rfcs/rfc3548.html

Below is the description of each bit represented in the PubChem Substructure Fingerprint. Some fingerprint bit descriptions are written in SMILES or SMARTS notation. For additional information on SMARTS and SMILES, please go to:

http://en.wikipedia.org/wiki/Simplified\_molecular\_input\_line\_entry\_sp
ecification

Section 1: Hierarchic Element Counts - These bits test for the presence or count of individual chemical atoms represented by their atomic symbol.

Bit Position	Bit Substructure
0	>= 4 H
1	>= 8 H
2	>= 16 H
3	>= 32 H
4	>= 1 Li
5	>= 2 Li
6	>= 1 B
7	>= 2 B
8	>= 4 B
9	>= 2 C
10	>= 4 C
11	>= 8 C
12	>= 16 C
13	>= 32 C
14	>= 1 N
15	>= 2 N
16	>= 4 N
17	>= 8 N
18	>= 1 0
19	>= 2 0
20	>= 4 0
21	>= 8 0
22	>= 16 0
23	>= 1 F
24	>= 2 F
25	>= 2 F >= 4 F
26	
27	
28	
29	A
30	>= 1 P
31	>= 2 P
32	>= 4 P
33	>= 1 S
34	>= 2 S
35	>= 4 S
36	>= 8 S
37	>= 1 Cl
38	>= 2 Cl
39	>= 4 Cl
40	>= 8 Cl
41	>= 1 K
42	>= 2 K
43	>= 1 Br
44	>= 2 Br
45	>= 4 Br
46	>= 1 I
47	>= 2 I
48	>= 4 I

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Section 1: Hierarchic Element Counts (cont.)

Bit Position	Bit Substructure
49	>= 1 Be
50	>= 1 Mg
51	
52	
53	>= 1 Sc
54	>= 1 Ti
55	>= 1 V
56	>= 1 Cr
57	>= 1 Mn
58	>= 1 Fe
59	>= 1 Co
60	>= 1 Ni
61	>= 1 Cu
62	>= 1 Zn
63	>= 1 Ga
64	>= 1 Ge
65	>= 1 As
66	>= 1 Se
67	>= 1 Kr
68	>= 1 Rb
69	>= 1 Sr
70	>= 1 Y
71	>= 1 Zr
72	>= 1 Nb
73	>= 1 Mo
74	>= 1 Ru
75	>= 1 Rh
76	>= 1 Pd
77	>= 1 Ag
78	>= 1 Cd
79	>= 1 In
80	>= 1 Sn
81	>= 1 Sb
82	>= 1 Te
83	>= 1 Xe
84	>= 1 Cs
85	>= 1 Ba
86	>= 1 Lu
87	>= 1 Hf
88	>= 1 Ta
89	>= 1 W
90	>= 1 Re
91	>= 1 Os
92	>= 1 Ir
93	>= 1 Pt
94	>= 1 Au
95	>= 1 Hg
96	>= 1 Tl
97	>= 1 Pb
98	>= 1 Bi

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Section 1: Hierarchic Element Counts (cont.)

Bit Position	Bit Substructure
99	>= 1 La
100	>= 1 Ce
101	>= 1 Pr
102	>= 1 Nd
103	>= 1 Pm
104	>= 1 Sm
105	>= 1 Eu
106	>= 1 Gd
107	>= 1 Tb
108	>= 1 Dy
109	>= 1 Ho
110	>= 1 Er
111	>= 1 Tm
112	>= 1 Yb
113	>= 1 Tc
114	>= 1 U

Section 2: Rings in a canonic Extended Smallest Set of Smallest Rings (ESSSR) ring set - These bits test for the presence or count of the described chemical ring system. An ESSSR ring is any ring which does not share three consecutive atoms with any other ring in the chemical structure. For example, naphthalene has three ESSSR rings (two phenyl fragments and the 10-membered envelope), while biphenyl will yield a count of only two ESSSR rings.

Bit Position	Bit Substructure
115	>= 1 any ring size 3
116	>= 1 saturated or aromatic carbon-only ring size 3
117	>= 1 saturated or aromatic nitrogen-containing ring size 3
118	>= 1 saturated or aromatic heteroatom-containing ring size 3
119	>= 1 unsaturated non-aromatic carbon-only ring size 3
120	>= 1 unsaturated non-aromatic nitrogen-containing ring size 3
121	>= 1 unsaturated non-aromatic heteroatom-containing ring size 3
122	>= 2 any ring size 3
123	>= 2 saturated or aromatic carbon-only ring size 3
124	>= 2 saturated or aromatic nitrogen-containing ring size 3
125	>= 2 saturated or aromatic heteroatom-containing ring size 3
126	>= 2 unsaturated non-aromatic carbon-only ring size 3
127	>= 2 unsaturated non-aromatic nitrogen-containing ring size 3
128	>= 2 unsaturated non-aromatic heteroatom-containing ring size 3
100	
129	>= 1 any ring size 4
130	>= 1 saturated or aromatic carbon-only ring size 4
131	>= 1 saturated or aromatic nitrogen-containing ring size 4
132	>= 1 saturated or aromatic heteroatom-containing ring size 4
133	>= 1 unsaturated non-aromatic carbon-only ring size 4

Section 2: Rings in a canonic ESSR ring set (cont.)

Bit Position	Bit Substructure
134	>= 1 unsaturated non-aromatic nitrogen-containing ring size 4
135	>= 1 unsaturated non-aromatic heteroatom-containing ring size 4
136	>= 2 any ring size 4
137	>= 2 saturated or aromatic carbon-only ring size 4
138	>= 2 saturated or aromatic nitrogen-containing ring size 4
139	>= 2 saturated or aromatic heteroatom-containing ring size 4
140	>= 2 unsaturated non-aromatic carbon-only ring size 4
141	>= 2 unsaturated non-aromatic nitrogen-containing ring size 4
142	>= 2 unsaturated non-aromatic heteroatom-containing ring size 4
143	>= 1 any ring size 5
144	>= 1 saturated or aromatic carbon-only ring size 5
145	>= 1 saturated or aromatic nitrogen-containing ring size 5
146	>= 1 saturated or aromatic heteroatom-containing ring size 5
147	>= 1 unsaturated non-aromatic carbon-only ring size 5
148	>= 1 unsaturated non-aromatic nitrogen-containing ring size 5
149	>= 1 unsaturated non-aromatic heteroatom-containing ring size 5
150	>= 2 any ring size 5
151	>= 2 saturated or aromatic carbon-only ring size 5
152	>= 2 saturated or aromatic nitrogen-containing ring size 5
153	>= 2 saturated or aromatic heteroatom-containing ring size 5
154	>= 2 unsaturated non-aromatic carbon-only ring size 5
155	>= 2 unsaturated non-aromatic nitrogen-containing ring size 5
156	>= 2 unsaturated non-aromatic heteroatom-containing ring size 5
157	>= 3 any ring size 5
158	>= 3 saturated or aromatic carbon-only ring size 5
159	>= 3 saturated or aromatic nitrogen-containing ring size 5
160	>= 3 saturated or aromatic heteroatom-containing ring size 5
161	>= 3 unsaturated non-aromatic carbon-only ring size 5
162	>= 3 unsaturated non-aromatic nitrogen-containing ring size 5
163	>= 3 unsaturated non-aromatic heteroatom-containing ring size 5
164	>= 4 any ring size 5
165	>= 4 saturated or aromatic carbon-only ring size 5
166	>= 4 saturated or aromatic nitrogen-containing ring size 5
167	>= 4 saturated or aromatic heteroatom-containing ring size 5
168	>= 4 unsaturated non-aromatic carbon-only ring size 5
169	>= 4 unsaturated non-aromatic nitrogen-containing ring size 5
170 171	>= 4 unsaturated non-aromatic heteroatom-containing ring size 5
172	<pre>&gt;= 5 any ring size 5 &gt;= 5 saturated or aromatic carbon-only ring size 5</pre>
173	>= 5 saturated or aromatic carbon-only fing size 5 >= 5 saturated or aromatic nitrogen-containing ring size 5
174	>= 5 saturated or aromatic heteroatom-containing ring size 5
175	>= 5 unsaturated non-aromatic carbon-only ring size 5
176	>= 5 unsaturated non-aromatic nitrogen-containing ring size 5
177	>= 5 unsaturated non-aromatic heteroatom-containing ring size 5 >= 5 unsaturated non-aromatic heteroatom-containing ring size 5
178	>= 1 any ring size 6
179	>= 1 saturated or aromatic carbon-only ring size 6
180	>= 1 saturated or aromatic nitrogen-containing ring size 6
181	>= 1 saturated or aromatic heteroatom-containing ring size 6
182	>= 1 unsaturated non-aromatic carbon-only ring size 6

Section 2: Rings in a canonic ESSR ring set (cont.)

Bit Position	Bit Substructure
183	>= 1 unsaturated non-aromatic nitrogen-containing ring size 6
184	>= 1 unsaturated non-aromatic heteroatom-containing ring size 6
185	>= 2 any ring size 6
186	>= 2 saturated or aromatic carbon-only ring size 6
187	>= 2 saturated or aromatic nitrogen-containing ring size 6
188	>= 2 saturated or aromatic heteroatom-containing ring size 6
189	>= 2 unsaturated non-aromatic carbon-only ring size 6
190	>= 2 unsaturated non-aromatic nitrogen-containing ring size 6
191	>= 2 unsaturated non-aromatic heteroatom-containing ring size 6
192	>= 3 any ring size 6
193	>= 3 saturated or aromatic carbon-only ring size 6
194	>= 3 saturated or aromatic nitrogen-containing ring size 6
195	>= 3 saturated or aromatic heteroatom-containing ring size 6
196	>= 3 unsaturated non-aromatic carbon-only ring size 6
197	>= 3 unsaturated non-aromatic nitrogen-containing ring size 6
198	>= 3 unsaturated non-aromatic heteroatom-containing ring size 6
199	>= 4 any ring size 6
200	>= 4 saturated or aromatic carbon-only ring size 6
201	>= 4 saturated or aromatic nitrogen-containing ring size 6
202	>= 4 saturated or aromatic heteroatom-containing ring size 6
203	>= 4 unsaturated non-aromatic carbon-only ring size 6
204	>= 4 unsaturated non-aromatic nitrogen-containing ring size 6
205	>= 4 unsaturated non-aromatic heteroatom-containing ring size 6
206	>= 5 any ring size 6
207	>= 5 saturated or aromatic carbon-only ring size 6
208	>= 5 saturated or aromatic nitrogen-containing ring size 6
209	>= 5 saturated or aromatic heteroatom-containing ring size 6
210	>= 5 unsaturated non-aromatic carbon-only ring size 6
211	>= 5 unsaturated non-aromatic nitrogen-containing ring size 6
212	>= 5 unsaturated non-aromatic heteroatom-containing ring size 6
01.2	>= 1 any ming gigo 7
213	>= 1 any ring size 7
214 215	>= 1 saturated or aromatic carbon-only ring size 7
216	<pre>&gt;= 1 saturated or aromatic nitrogen-containing ring size 7 &gt;= 1 saturated or aromatic heteroatom-containing ring size 7</pre>
217	>= 1 unsaturated non-aromatic carbon-only ring size 7
217	>= 1 unsaturated non-aromatic nitrogen-containing ring size 7
219	>= 1 unsaturated non-aromatic heteroatom-containing ring size 7
220	>= 2 any ring size 7
221	>= 2 saturated or aromatic carbon-only ring size 7
222	>= 2 saturated or aromatic nitrogen-containing ring size 7
223	>= 2 saturated or aromatic heteroatom-containing ring size 7
224	>= 2 unsaturated non-aromatic carbon-only ring size 7
225	>= 2 unsaturated non-aromatic carbon-only ring size / >= 2 unsaturated non-aromatic nitrogen-containing ring size 7
226	>= 2 unsaturated non-aromatic heteroatom-containing ring size 7
227	>= 1 any ring size 8
228	>= 1 saturated or aromatic carbon-only ring size 8
229	>= 1 saturated or aromatic nitrogen-containing ring size 8
230	>= 1 saturated or aromatic heteroatom-containing ring size 8
231	>= 1 unsaturated non-aromatic carbon-only ring size 8

Section 2: Rings in a canonic ESSR ring set (cont.)

Bit Position	Bit Substructure
232	>= 1 unsaturated non-aromatic nitrogen-containing ring size 8
233	>= 1 unsaturated non-aromatic heteroatom-containing ring size 8
234	>= 2 any ring size 8
235	>= 2 saturated or aromatic carbon-only ring size 8
236	>= 2 saturated or aromatic nitrogen-containing ring size 8
237	>= 2 saturated or aromatic heteroatom-containing ring size 8
238	>= 2 unsaturated non-aromatic carbon-only ring size 8
239	>= 2 unsaturated non-aromatic nitrogen-containing ring size 8
240	>= 2 unsaturated non-aromatic heteroatom-containing ring size 8
241	>= 1 any ring size 9
242	>= 1 saturated or aromatic carbon-only ring size 9
243	>= 1 saturated or aromatic nitrogen-containing ring size 9
244	>= 1 saturated or aromatic heteroatom-containing ring size 9
245	>= 1 unsaturated non-aromatic carbon-only ring size 9
246	>= 1 unsaturated non-aromatic nitrogen-containing ring size 9
247	>= 1 unsaturated non-aromatic heteroatom-containing ring size 9
248	>= 1 any ring size 10
249	>= 1 saturated or aromatic carbon-only ring size 10
250	>= 1 saturated or aromatic nitrogen-containing ring size 10
251	>= 1 saturated or aromatic heteroatom-containing ring size 10
252	>= 1 unsaturated non-aromatic carbon-only ring size 10
253	>= 1 unsaturated non-aromatic nitrogen-containing ring size 10
254	>= 1 unsaturated non-aromatic heteroatom-containing ring size 10
255	>= 1 aromatic ring
256	>= 1 hetero-aromatic ring
257	>= 2 aromatic rings
258	>= 2 hetero-aromatic rings
259	>= 3 aromatic rings
260	>= 3 hetero-aromatic rings
261	>= 4 aromatic rings
262	>= 4 hetero-aromatic rings

**Section 3:** Simple atom pairs - These bits test for the presence of patterns of bonded atom pairs, regardless of bond order or count.

Bit Position	Bit Substructure
263	Li-H
264	Li-Li
265	Li-B
266	Li-C
267	Li-O
268	Li-F
269	Li-P
270	Li-S

Section 3: Simple atom pairs (cont.)

Bit Position	Bit Substructure
271	Li-Cl
272	В-Н
273	B-B
274	B-C
275	B-N
276	B-O
277	B-F
278	B-Si
279	B-P
280	B-S
281	B-Cl
282	B-Br
283	C-H
284	C-C
285	C-N
286	C-0
287	C-F
288	C-Na
289	C-Mg
290	C-Al
291	C-Si
292	C-P
293	C-S
294	C-Cl
295	C-As
296	C-Se
297	C-Br
298	C-I
299	N-H
300	N-N
301	N-O
302	N-F
303	N-Si
304	N-P
305	N-S
306	N-Cl
307 308	N-Br
	O-H O-O
309	
310 311	O-Mg
312	O-Na
313	O-Al O-Si
313	0-S1 0-P
314	0-P 0-K
316	0-K F-P
317	F-S
318	Al-H
319	Al-Cl
320	Si-H
320	PT-U

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Section 3: Simple atom pairs (cont.)

Bit Position	Bit Substructure
321	Si-Si
322	Si-Cl
323	P-H
324	P-P
325	As-H
326	As-As

Section 4: Simple atom nearest neighbors - These bits test for the presence of atom nearest neighbor patterns, regardless of bond order (denoted by "~") or count, but where bond aromaticity (denoted by ":") is significant.

Bit Position	Bit Substructure
327	C(~Br)(~C)
328	C(~Br)(~C)(~C)
329	C(~Br)(~H)
330	C(~Br)(:C)
331	C(~Br)(:N)
332	C(~C)(~C)
333	C(~C)(~C)(~C)
334	C(~C)(~C)(~C)(~C)
335	C(~C)(~C)(~C)(~H)
336	C(~C)(~C)(~C)(~N)
337	C(~C)(~C)(~C)(~O)
338	C(~C)(~C)(~H)(~N)
339	C(~C)(~C)(~H)(~O)
340	C(~C)(~C)(~N)
341	C(~C)(~C)(~O)
342	C(~C)(~Cl)
343	C(~C)(~Cl)(~H)
344	C(~C)(~H)
345	C(~C)(~H)(~N)
346	C(~C)(~H)(~O)
347	C(~C)(~H)(~O)(~O)
348	C(~C)(~H)(~P)
349	C(~C)(~H)(~S)
350	C(~C)(~I)
351	C(~C)(~N)
352	C(~C)(~O)
353	C(~C)(~S)
354	C(~C)(~Si)
355	C(~C)(:C)
356	C(~C)(:C)(:C)
357	$C(\sim C)(:C)(:N)$
358	C(~C)(:N)
359	$C(\sim C)(:N)(:N)$
360	C(~Cl)(~Cl)

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**Section 4:** Simple atom nearest neighbors (cont.)

Bit Position	Bit Substructure
361	C(~Cl)(~H)
362	C(~Cl)(:C)
363	C(~F)(~F)
364	C(~F)(:C)
365	C(~H)(~N)
366	C(~H)(~O)
367	C(~H)(~O)(~O)
368	C(~H)(~S)
369 370	C(~H)(~Si) C(~H)(:C)
370	C(~H)(:C)(:C)
372	$C(\sim H)(:C)(:C)$
373	C(~H)(:N)
374	C(~H)(~H)(~H)
375	C(~N)(~N)
376	$C(\sim N)(:C)$
377	$C(\sim N)(:C)(:C)$
378	$C(\sim N)(:C)(:N)$
379	$C(\sim N)(:N)$
380	C(~O)(~O)
381	C(~O)(:C)
382	C(~O)(:C)(:C)
383	C(~S)(:C)
384 385	C(:C)(:C) C(:C)(:C)(:C)
386	C(:C)(:C)(:C)
387	C(:C)(:N)
388	C(:C)(:N)(:N)
389	C(:N)(:N)
390	N(~C)(~C)
391	N(~C)(~C)(~C)
392	N(~C)(~C)(~H)
393	$N(\sim C)(\sim H)$
394	N(~C)(~H)(~N)
395	N(~C)(~O)
396	N(~C)(:C)
397	$N(\sim C)(:C)(:C)$
398	N(~H)(~N)
399 400	N(~H)(:C) N(~H)(:C)(:C)
401	N(~H)(.C)(.C) N(~O)(~O)
402	N(~O)(:O)
403	N(:C)(:C)
404	N(:C)(:C)(:C)
405	O(~C)(~C)
406	O(~C)(~H)
407	O(~C)(~P)
408	O(~H)(~S)
409	O(:C)(:C)
410	P(~C)(~C)

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#### PubChem Substructure Fingerprint Description (cont.)

**Section 4:** Simple atom nearest neighbors (cont.)

Bit Position	Bit Substructure
411	P(~O)(~O)
412	S(~C)(~C)
413	S(~C)(~H)
414	S(~C)(~O)
415	Si(~C)(~C)

Section 5: Detailed atom neighborhoods - These bits test for the
 presence of detailed atom neighborhood patterns, regardless
 of count, but where bond orders are specific, bond
 aromaticity matches both single and double bonds, and where
 "-", "=", and "#" matches a single bond, double bond, and
 triple bond order, respectively.

Bit Position	Bit Substructure
416	C=C
417	C#C
418	C=N
419	C#N
420	C=0
421	C=S
422	N=N
423	N=O
424	N=P
425	P=0
426	P=P
427	C(#C)(-C)
428	C(#C)(-H)
429	C(#N)(-C)
430	C(-C)(-C)(=C)
431	C(-C)(-C)(=N)
432	C(-C)(-C)(=0)
433	C(-C)(-C1)(=0)
434	C(-C)(-H)(=C)
435	C(-C)(-H)(=N)
436	C(-C)(-H)(=0)
437	C(-C)(-N)(=C)
438	C(-C)(-N)(=N)
439	C(-C)(-N)(=0)
440	C(-C)(-O)(=O)
441	C(-C)(=C)
442	C(-C)(=N)
443	C(-C)(=0)
444	C(-Cl)(=0)
445	C(-H)(-N)(=C)
446	C(-H)(=C)
447	C(-H)(=N)
448	C(-H)(=O)

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**Section 5:** Detailed atom neighborhoods (cont.)

Bit Position	Bit Substructure
449	C(-N)(=C)
450	C(-N)(=N)
451	C(-N)(=O)
452	C(-O)(=O)
453	N(-C)(=C)
454	N(-C)(=O)
455	N(-O)(=O)
456	P(-O)(=O)
457	S(-C)(=O)
458	S(-O)(=O)
459	S(=0)(=0)

**Section 6:** Simple SMARTS patterns - These bits test for the presence of simple SMARTS patterns, regardless of count, but where bond orders are specific and bond aromaticity matches both single and double bonds.

Die Desieles	Dit Colombon
Bit Position 460	Bit Substructure
	C-C-C#C
461	O-C-C=N
462	0-C-C=0
463	N:C-S-[#1]
464	N-C-C=C
465	0=S-C-C
466	N#C-C=C
467	C=N-N-C
468	O=S-C-N
469	S-S-C:C
470	C:C-C=C
471	S:C:C:C
472	C:N:C-C
473	S-C:N:C
474	S:C:C:N
475	S-C=N-C
476	C-O-C=C
477	N-N-C:C
478	S-C=N-[#1]
479	S-C-S-C
480	C:S:C-C
481	O-S-C:C
482	C:N-C:C
483	N-S-C:C
484	N-C:N:C
485	N:C:C:N
486	N-C:N:N
487	N-C=N-C
488	N-C=N-[#1]

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Section 6: Simple SMARTS patterns (cont.)

Bit Position	Bit Substructure
489	N-C-S-C
490	C-C-C=C
491	C-N:C-[#1]
492	N-C:O:C
493	O=C-C:C
494	O=C-C:N
495	C-N-C:C
496	N:N-C-[#1]
497	O-C:C:N
498	O-C=C-C
499	N-C:C:N
500	C-S-C:C
501	Cl-C:C-C
502	N-C=C-[#1]
503	Cl-C:C-[#1]
504	N:C:N-C
505	C1-C:C-O
506	C-C:N:C
507	C-C-S-C
508	S=C-N-C
509	Br-C:C-C
510	[#1]-N-N-[#1]
511	S=C-N-[#1]
512	C-[As]-O-[#1]
513	S:C:C-[#1]
514	O-N-C-C
515	N-N-C-C
516	[#1]-C=C-[#1]
517	N-N-C-N
518	O=C-N-N
519	N=C-N-C
520	C=C-C:C
521	C:N-C-[#1]
522	C-N-N-[#1]
523	N:C:C-C
524	C-C=C-C
525	[As]-C:C-[#1]
526	Cl-C:C-Cl
527	C:C:N-[#1]
528	[#1]-N-C-[#1]
529	Cl-C-C-Cl
530	N:C-C:C
531	S-C:C-C
532	S-C:C-[#1]
533	S-C:C-N
534	S-C:C-0
535	O=C-C-C
536	O=C-C-N
537	O=C-C-O
538	N=C-C-C

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Section 6: Simple SMARTS patterns (cont.)

Bit Position	Bit Substructure
539	N=C-C-[#1]
540	C-N-C-[#1]
541	O-C:C-C
542	O-C:C-[#1]
543	O-C:C-N
544	0-C:C-N
545	N-C:C-C
546	N-C:C-[#1]
547	N-C:C-N
548	O-C-C:C
549	N-C-C:C
550	Cl-C-C-C
551	C1-C-C-O
552	C:C-C:C
553	O=C-C=C
554	Br-C-C-C
555	N=C-C=C
556	C=C-C-C
557	N:C-O-[#1]
558	O=N-C:C
559	O-C-N-[#1]
560	N-C-N-C
561	Cl-C-C=O
562	Br-C-C=O
563	O-C-O-C
564	C=C-C=C
565	C:C-O-C
566	O-C-C-N
567	O-C-C-O
568	N#C-C-C
569	N-C-C-N
570	C:C-C-C
571	[#1]-C-O-[#1]
572	N:C:N:C
573	O-C-C=C
574	0-C-C:C-C
575	0-C-C:C-0
576	N=C-C:C-[#1]
577	C:C-N-C:C
578	C-C:C-C:C
579	O=C-C-C
580	O=C-C-C-N
581	O=C-C-C-O
582	C-C-C-C-C
583	Cl-C:C-O-C
584	C:C-C=C-C
585	C-C:C-N-C
586	C-S-C-C-C
587	N-C:C-O-[#1]
588	0=C-C-C=0
500	0-0 0 0-0

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Section 6: Simple SMARTS patterns (cont.)

Bit Position	Bit Substructure
	C-C:C-O-C
589	
590	C-C:C-O-[#1]
591	Cl-C-C-C
592	N-C-C-C
593	N-C-C-C-N
594	C-O-C-C=C
595	C:C-C-C-C
596	N=C-N-C-C
597	O=C-C-C:C
598	C1-C:C:C-C
599	[#1]-C-C=C-[#1]
600	N-C:C:C-C
601	N-C:C:C-N
602	O=C-C-N-C
603	C-C:C:C-C
604	C-O-C-C:C
605	O=C-C-O-C
606	O-C:C-C-C
607	N-C-C-C:C
608	C-C-C-C:C
609	Cl-C-C-N-C
610	C-O-C-O-C
611	N-C-C-N-C
612	N-C-O-C-C
613	C-N-C-C-C
614	C-C-O-C-C
615	N-C-C-O-C
616	C:C:N:N:C
617	C-C-C-O-[#1]
618	
	C:C-C-C:C
619	0-C-C=C-C
620	C:C-O-C-C
621	N-C:C:C:N
622	O=C-O-C:C
623	O=C-C:C-C
624	O=C-C:C-N
625	O=C-C:C-O
626	C-O-C:C-C
627	O=[As]-C:C:C
628	C-N-C-C:C
629	S-C:C:C-N
630	O-C:C-O-C
631	O-C:C-O-[#1]
632	C-C-O-C:C
633	N-C-C:C-C
634	C-C-C:C-C
635	N-N-C-N-[#1]
636	= :: =
637	C-N-C-N-C
	0-C-C-C-C
638	O-C-C-C-N

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Section 6: Simple SMARTS patterns (cont.)

Bit Position	Bit Substructure
639	0-C-C-C-O
640	C=C-C-C-C
641	O-C-C-C=C
642	O-C-C-C=O
643	[#1]-C-C-N-[#1]
644	C-C=N-N-C
645	O=C-N-C-C
646	O=C-N-C-[#1]
647	O=C-N-C-N
648	O=N-C:C-N
649	O=N-C:C-O
650	O=C-N-C=O
651	O-C:C:C-C
652	O-C:C:C-N
653	0-C:C:C-O
654	N-C-N-C-C
655	O-C-C-C:C
656	C-C-N-C-C
657	C-N-C:C-C
658	C-N-C-C-C C-C-S-C-C
659	O-C-C-N-C
660	C-C=C-C-C
661	0-C-O-C-C
662	0-C-C-O-C
663	O-C-C-O-[#1]
664	C-C=C-C=C
665	N-C:C-C-C
666	C=C-C-O-C
667	C=C-C-O-[#1]
668	C-C:C-C-C
669	C1-C:C-C=O
670	Br-C:C:C-C
671	O=C-C=C-C
672	O=C-C=C-[#1]
673	O=C-C=C-N
674	N-C-N-C:C
675	Br-C-C-C:C
676	N#C-C-C-C
677	C-C=C-C:C
678	C-C-C=C-C
679	C-C-C-C-C-C
680	O-C-C-C-C
681	O-C-C-C-C
682	O-C-C-C-N
683	N-C-C-C-C
684	O=C-C-C-C
685	O=C-C-C-C-N
686	0=C-C-C-C-O
687	O=C-C-C-C=O
688	C-C-C-C-C-C

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**Section 6:** Simple SMARTS patterns (cont.)

Bit Position	Bit Substructure
689	O-C-C-C-C-C
690	O-C-C-C-C-O
691	O-C-C-C-C-C-N
692	O=C-C-C-C-C
693	O=C-C-C-C-O
694	O=C-C-C-C-C=O
695	O=C-C-C-C-C-N
696	C-C-C-C-C-C-C
697	C-C-C-C-C-C(C)-C
698	O-C-C-C-C-C-C
699	O-C-C-C-C(C)-C
700	O-C-C-C-C-C-O-C
701	O-C-C-C-C(O)-C
702	O-C-C-C-C-C-N-C
703	O-C-C-C-C-C(N)-C
704	O=C-C-C-C-C-C
705	O=C-C-C-C(O)-C
706	O=C-C-C-C(=O)-C
707	O=C-C-C-C-C(N)-C
708	C-C(C)-C-C
709	C-C(C)-C-C-C
710	C-C-C(C)-C-C
711	C-C(C)(C)-C-C
712	C-C(C)-C(C)-C

**Section 7:** Complex SMARTS patterns - These bits test for the presence of complex SMARTS patterns, regardless of count, but where bond orders and bond aromaticity are specific.

Bit Position	<u>Bit Substructure</u>
713	Cclccc(C)ccl
714	Cclccc(O)ccl
715	Cclccc(S)ccl
716	Cclccc(N)ccl
717	Cc1ccc(Cl)cc1
718	Cclccc(Br)ccl
719	Oc1ccc(O)cc1
720	Oclccc(S)ccl
721	Oclccc(N)ccl
722	Oclccc(Cl)ccl
723	Oclccc(Br)ccl
724	Sclccc(S)ccl
725	Sclccc(N)ccl
726	Sc1ccc(Cl)cc1
727	Sclccc(Br)ccl
728	Nc1ccc(N)cc1
729	Nc1ccc(Cl)cc1
730	Nclccc(Br)ccl

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Section 7: Complex SMARTS patterns (cont.)

Bit Position	Bit Substructure
731	Clc1ccc(Cl)cc1
732	Clc1ccc(Br)cc1
733	Brclccc(Br)ccl
734	Cclcc(C)cccl
735	Cclcc(O)cccl
736	Cclcc(S)cccl
737	Cclcc(N)cccl
738	Cclcc(Cl)cccl
739	Cclcc(Br)cccl
740	Oclcc(O)cccl
741	Oclcc(S)cccl
742	Oclcc(N)cccl
743	Oclcc(Cl)cccl
744	
745	Oclcc(Br)cccl
	Sclcc(S)cccl
746	Sclcc(N)cccl
747	Sclcc(Cl)cccl
748	Sclcc(Br)cccl
749	Nclcc(N)cccl
750	Nclcc(Cl)cccl
751	Nclcc(Br)cccl
752	Clclcc(Cl)cccl
753	Clclcc(Br)cccl
754	Brclcc(Br)cccl
755	Cclc(C)ccccl
756	Cclc(O)ccccl
757	Cclc(S)ccccl
758	Cclc(N)ccccl
759	Cclc(Cl)ccccl
760	Cclc(Br)ccccl
761	Oc1c(O)cccc1
762	Oclc(S)ccccl
763	Oclc(N)ccccl
764	Oclc(Cl)ccccl
765	Oclc(Br)ccccl
766	Sc1c(S)cccc1
767	Sc1c(N)cccc1
768	Sc1c(Cl)cccc1
769	Sclc(Br)ccccl
770	Nc1c(N)cccc1
771	Nc1c(Cl)cccc1
772	Nclc(Br)ccccl
773	Clc1c(Cl)cccc1
774	Clc1c(Br)cccc1
775	Brc1c(Br)cccc1
776	CC1CCC(C)CC1
777	CC1CCC(O)CC1
778	CC1CCC(S)CC1
779	CC1CCC(N)CC1
780	CC1CCC(C1)CC1
700	CCTCCC(CT)CCT

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Section 7: Complex SMARTS patterns (cont.)

Bit Position	Bit Substructure
781	CC1CCC(Br)CC1
782	OC1CCC(O)CC1
783	OC1CCC(S)CC1
784	OC1CCC(N)CC1
785	OC1CCC(C1)CC1
786	OC1CCC(Br)CC1
787	SC1CCC(S)CC1
788	SC1CCC(N)CC1
789	SC1CCC(C1)CC1
790	
	SC1CCC(Br)CC1
791	NC1CCC(N)CC1
792	NC1CCC(C1)CC1
793	NC1CCC(Br)CC1
794	ClC1CCC(Cl)CC1
795	ClC1CCC(Br)CC1
796	BrC1CCC(Br)CC1
797	CC1CC(C)CCC1
798	CC1CC(O)CCC1
799	CC1CC(S)CCC1
800	CC1CC(N)CCC1
801	CC1CC(C1)CCC1
802	CC1CC(Br)CCC1
803	OC1CC(O)CCC1
804	OC1CC(S)CCC1
	OCICC(S)CCCI
805	
806	OC1CC(C1)CCC1
807	OC1CC(Br)CCC1
808	SC1CC(S)CCC1
809	SC1CC(N)CCC1
810	SC1CC(C1)CCC1
811	SC1CC(Br)CCC1
812	NC1CC(N)CCC1
813	NC1CC(Cl)CCC1
814	NC1CC(Br)CCC1
815	ClC1CC(Cl)CCC1
816	ClC1CC(Br)CCC1
817	BrC1CC(Br)CCC1
818	CC1C(C)CCCC1
819	CC1C(0)CCCC1
820	CC1C(S)CCCC1
821	CC1C(N)CCCC1
822	CC1C(C1)CCCC1
823	CC1C(Br)CCCC1
824	
	0C1C(0)CCCC1
825	OC1C(S)CCCC1
826	OC1C(N)CCCC1
827	OC1C(C1)CCCC1
828	OC1C(Br)CCCC1
829	SC1C(S)CCCC1
830	SC1C(N)CCCC1

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Section 7: Complex SMARTS patterns (cont.)

Bit Position	Bit Substructure
831	SC1C(C1)CCCC1
832	SC1C(Br)CCCC1
833	NC1C(N)CCCC1
834	NC1C(Cl)CCCC1
835	NC1C(Br)CCCC1
836	clc1c(cl)cccc1
837	ClC1C(Br)CCCC1
838	BrC1C(Br)CCCC1
839	CC1CC(C)CC1
840	CC1CC(O)CC1
841	CC1CC(S)CC1
842	CC1CC(N)CC1
843	CC1CC(N)CC1
844	CC1CC(Br)CC1
845	
	OC1CC(0)CC1
846	OC1CC(S)CC1
847	OC1CC(N)CC1
848	OC1CC(C1)CC1
849	OC1CC(Br)CC1
850	SC1CC(S)CC1
851	SC1CC(N)CC1
852	SC1CC(C1)CC1
853	SC1CC(Br)CC1
854	NC1CC(N)CC1
855	NC1CC(C1)CC1
856	NC1CC(Br)CC1
857	ClC1CC(Cl)CC1
858	ClC1CC(Br)CC1
859	BrC1CC(Br)CC1
860	CC1C(C)CCC1
861	CC1C(O)CCC1
862	CC1C(S)CCC1
863	CC1C(N)CCC1
864	CC1C(Cl)CCC1
865	CC1C(Br)CCC1
866	OC1C(O)CCC1
867	OC1C(S)CCC1
868	OC1C(N)CCC1
869	OC1C(C1)CCC1
870	OC1C(Br)CCC1
871	SC1C(S)CCC1
872	SC1C(N)CCC1
873	SC1C(C1)CCC1
874	SC1C(Br)CCC1
875	NC1C(N)CCC1
876	NC1C(C1)CC1
877	NC1C(Br)CCC1
878	ClC1C(Cl)CCC1
879	ClC1C(Br)CCC1
880	BrC1C(Br)CCC1

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#### Decoding PubChem Fingerprints

PubChem fingerprints are currently 881 bits in length. Binary data is stored in one byte increments. The fingerprint is, therefore, 111 bytes in length (888 bits), which includes padding of seven bits at the end to complete the last byte. A four-byte prefix, containing the bit length of the fingerprint (881 bits), increases the stored PubChem fingerprint size to 115 bytes (920 bits).

When PubChem fingerprints are encoded in base64 format, the base64-encoded fingerprints are 156 bytes in length. The last two bytes are padding so that the base64 length is divisible by four (156 bytes - 2 bytes = 154 bytes). Each base64 byte encodes six binary bits (154 bytes \* 6 bits/byte = 924 bits). The last four bits are padding to complete the last base64 byte (924 bits - 4 bits = 920 bits). The resulting 920 binary bits (115 bytes) are described in the previous paragraph.

#### Document Version History

- V1.3 2009May01 Updated introduction to describe how to identify the PubChem Substructure Fingerprint property in a PubChem Compound record.
- V1.2 2007Aug30 Added section on decoding PubChem fingerprints.
- V1.1 2007Aug06 Corrected and expanded documentation of bits with SMARTS patterns used.
- V1.0 2005Dec02 Initial release.