## **Generic Searching by Use of Rotated Formula Indexes\***

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The term, "generic searching," is variously interpreted to mean classified, non-specific, group or categorized searching. For research chemists, generic searches are a valuable tool. Structurally related chemicals with a specified combination of functional groups, rings, or other desiderata used to code or describe chemicals, may be found in a more orderly fashion by generic searching. While many machine searching systems code molecular formulas, this information is rarely used in generic searches. The Chemical Biological Coordination Center (CBCC) file provided molecular formulas which were not used much in searching.

Generic searching often is done on a less sophisticated level. For example, one researcher regularly looks for articles on new uses of lithium. Obviously, as a lithium manufacturer, his company is interested in finding new applications. Similarly, chemists in the missile propellant field are interested in all new boron compounds since many are valuable in the production of propellants.

From my own experience, I have found that the conventional formula indexes of *Chemical Abstracts* and of *Index Chemicus* do not permit a convenient means of searching generically except for certain classes such as steroids (C17), etc. For example, a search for "all compounds in which phosphorus is attached to at least five oxygen atoms" is completely impossible unless you search every page of the index. If L. Pauling and V. Schomaker had been able to make such a generic search, it might have avoided the need to retract statements made in a communication commenting upon the works of E. Ronwin.

In November, 1951, Ronwin published a paper in J. Am. Chem. Soc.<sup>1</sup> introducing a new structural formula for both ribo- and desoxyribonucleic acids. The structure had as its core a  $(P_2O_5)_n$  polymer chain of phosphoanhydride links. By a theoretical treatment, Ronwin showed the formula to be compatible with available factual data concerned with the structure of nucleic acids. In February, 1952, L. Pauling and V. Schomaker wrote in a communication to the editor<sup>2</sup>: "In the proposed structure for the nucleic acids each P atom has five O atoms attached to it, three of which bind it to adjacent P atoms, and two of which are in a OH group and a sugar ester group, respectively. There is, however, no precedent for a structure in which P is bonded to five O atoms... the ligation of five O atoms about each P atom is such an unlikely structural feature that the proposed phosphotri-anhydride formula for the nucleic acids deserves no serious consideration."

However, in the July, 1952, issue,<sup>8</sup> they made a most gentlemanly retraction in a second communication: "Dr. Ronwin has now kindly informed us that he has become aware of earlier references in the literature to compounds to which structures have been attributed involving quinquepositive P bonded to five O atoms or to a total of five O atoms and similar atoms. Anschütz<sup>4</sup> prepared four compounds to which he assigned structures involving ligation of five O atoms to a P atom ... our statement that there is no precedent for a structure in which a P atom is bonded to five O atoms must accordingly be withdrawn."

Frankly I can't recall how I even came across this interesting discussion. I was doing a search on a completely unrelated matter. It occurred to me at that time that the retraction would have been unnecessary—if a generic formula index to *Chemical Abstracts* were available. I was intrigued to learn how Dr. Ronwin managed to find the reference to Anschütz's paper and others mentioned by Pauling. It had been impossible for me to find them in a search. I wrote to Dr. Ronwin<sup>5</sup> and asked him how he had managed to find the necessary precedent. Dr. Ronwin replied<sup>6</sup> that a colleague had remembered the paper as an abstractor for *Chemisches Zentralblatt*!

In a subsequent letter, however, Prof. Pauling pointed out that the retraction did not involve a retraction of the criticism of the proposed phospho-tri-anhydride formulas for the nucleic acids. The retraction was limited to the statement that "there is no precedent for a structure in which a P atom is bonded to five O atoms." Prof. Pauling correctly states that there might be some question about the correctness of the structures assigned to some of these compounds, and especially that these compounds are extremely sensitive to moisture.<sup>7</sup>

This, then, is the explanation of why and when I first felt the need for a rotated formula index. I wonder to this day whether there might not have been other examples in the literature which Pauling or Ronwin might have found if a rotated index were available. This example illustrates very well the role that new scientific information services can play in advancing scientific theory. In a short paper such as this, there is little need to belabor the question of the potential value of generic indexes. However, generic formula indexes may not be sufficiently appreciated as a searching tool since they have not, until now, been made available.

<sup>\*</sup> Presented at the 141st National Meeting, American Chemical Society, Divison of Chemical Literature, March 22, 1962, Washington, D. C.

In previous papers, Skolnik<sup>8</sup> and Fletcher<sup>9</sup> have indicated that conventional formula indexes leave much to be desired. They have shown how an "inverted" filing system with carbon and hydrogen given the lowest filing priority can be useful, not only in shortening the search for a specific compound, but also in locating generically related compounds. In the Skolnik system used at the Hercules Powder Co., elements are filed alphabetically with carbon at the end of the formula and hydrogen completely ignored. In this system structural formulas are available on the file cards. At American Cyanamid, the Fletcher system, based on the periodic order of the elements, is used. It has many definite advantages for grouping together structurally related chemicals. Dyson<sup>10</sup> also used a similar method. However, in any of these systems, there is only one file assignment made for each chemical. In order to find a particular compound, it is necessary to anticipate all possible positions where a particular element might occur. For example, if a chemical contains both aluminum and phosphorus, and the chemist is searching for phosphorus compounds, he may not find it under P, but rather under Al.

In previous papers that I have presented before this Division, I discussed the methods employed in preparing the molecular formula indexes to *Index Chemicus*.<sup>11, 12</sup> Briefly, this involves the preparation of a single punchedcard for each formula. The cards are converted to magnetic tape and sorted on an electronic computer and printed on a high speed printer for photo-offset reproduction. Recently, over 150,000 formulas were processed in the preparation of our first two-year cumulation. In this cumulation, we completely redesigned the index format to make it easier to use as well as to save space and lower costs. As a completely by-product operation, we have also prepared the first rotated formula index, a *RotaForm Index*. Let me first describe the *RotaForm Index* and then let me illustrate its use with several examples.

In Fig. 1, you see a typical page from the old cumulative molecular formula index to the Index Chemicus. As in Chemical Abstracts, the Hill system giving priority to carbon and hydrogen is used. All subsequent elements are filed alphabetically. However, a new heading has been created for each carbon-hydrogen combination to speed up locating a desired compound, simultaneously reducing the size of the index by about 25%. This can be seen by examining Fig. 2 which shows the new format. Notice the considerable amount of wasted white space below the carbon-hydrogen headings in the old format. The second major change in the format is the way that the serial numbers of the individual compounds are listed. Whereas each address was listed on one line on the left, they are now on the right of the molecular formula and one or more serial numbers appear after a dotted line. Notice in most cases that the line containing a new molecular formula is always justified on both margins.

As I mentioned before, one card is prepared for each formula in the molecular formula index used now. To prepare the *RotaForm Index*, the computer essentially duplicates the card as many times as there are different elements. If the compound contains, in addition to C and H, five other different elements, the formula will be repeated in the index five times. Thus, C23 H20 Al2 Br3 F4 Na2 P3 (hypothetical example) would be repeated in the *RotaForm Index* five times, once each under the following arrangements:

- (1) Al2 Br3 F4 Na2 P3 C23 H2O
- (2) Br3 Al2 F4 Na2 P3 C23 H2O
  (3) F4 Al2 Br3 Na2 P3 O23 H2O
- (4) Na2 Al2 Br3 F4 P3 C23 H2O
- (f)  $P_3 Al2 Br_3 F_4 N_{a2} C_{23} H_{20}$

With this type of index, a search for compounds containing any one of these elements is possible in a quick and orderly fashion. Fig. 3 is a page from the new *RotaForm Index*. This page contains all the formulas containing three phosphorus atoms. Since all of the elements are sorted alphabetically, it is a simple matter to locate, for example, all compounds containing three phosphorus atoms and also containing fluorine. These are found by scanning all of the P3 compound listings until P3 Fl2. There cannot be any fluorine compounds beyond this group because of the alphabetical arrangement. Seven such compounds have been checked. A similar search for P4 compounds were located.

There are some obvious advantages of the RotaForm Index. It will now be quite easy to determine, for example, all reported compounds containing a particular element such as boron. Much more sophisticated generic searches can also be done, as shown in Fig. 4. (Dr. Joe Clark of Lederle Laboratories kindly supplied the questions and the searches were conducted by members of the Index Chemicus staff: Mrs. Eleanore H. Peitsch Baus, Assistant Managing Editor, and George F. Corkery, Indexer, and Miss Sandra Goldman, Secretary).

The RotaForm Index is a supplementary tool to the conventional formula index. It permits a limited range of generic searches and is particularly valuable for a search involving the less-frequently occurring elements. It is somewhat more cumbersome for searches involving O, N, and S. Even these, however, can be performed with amazing results considering the low cost involved. The RotaForm Index is not a panacea for the problems of generic searching. It is, however, one more tool to add to the chemist's armamentarium. It is a by-product of our regular formula indexing, and completely machine-made. It simply advises the reader which molecular formulas contain a particular combination of elements. In order to keep the size down, we have not repeated serial numbers, which can be found in the molecular formula index.

The searcher will save considerable time if he arranges molecular formulas found in the *RotaForm Index* by the number of carbon atoms before looking up the serial numbers. This is strictly a clerical procedure. In fact, any use of the *RotaForm Index* can be assigned to a clerk until the final screening of the structural diagrams.

Keep in mind that the use of this system in the average company chemical file, as is the case in the Hercules and in the American Cyanamid System, each molecular formula card is accompanied by the structural diagram. Such a card filing system eliminates the chore of going from the compound serial number to a register showing the structural diagram itself.

## ORGANIC COMPOUNDS

С,			$C_3H_2$		$C_4H_4$
6586- SC 1	AG F3 S	7736- 2C 2H 3 ALZ 17 0	4821- 4C 3M 2 BR CL3 F	2 6970-2 N3	5352-17C 4H 4 AG F3 04
7252- 3	8 F6 K	7868- 2 B CL2	7501-1 CL N D	4804-7 N3 04	5352-14 BR F3 0
6819- 1	CL2 FNO	7736-1 B I4 O	4802-3 CL3 N3 O	2 5530-2 03 P	7141-1 BR N S
6819- 3	CL2 F2 02 5	6864-1 BR CL D	4779-8 53	5824-1 03 P	6364- 3 BR N3 0
7207- 1	CL3 F 02 S	6335- 4 BR CL N 02	7892- 3C 3M 3 B3 N6	6063-10 04 P 5I	7501- 4 CL N 02
5464- 5	04 CL2 51	7356-1 CL3 0 5	7614-20 CL4 N 0	5524- 2 B CL4 N	5695-1 CL2 HG 04
5464- 3	D4 F2 SI	7356-2 F3 O S	7303-58 F3 N2 04	S 7041-1 B F4 N	7614-17 CL3 N 02
5023- 6	F2 0	5803-10 K 0 52	7356-3 F5 0 S	7041-13 CL N 04 7041-16 F6 N P	7599-7 CL6 S2 5626-3 CU2 12 N2
6586- 2	F2 5	4705- 2 N3 52	4901-3 1 52	6135-1 N3 04 P	6378- 2 D2 04
7049- 1	MN N3 04	5859-68 N4 NA S	7769-1 N 02 S	4858-44 NA P SZ	5352-2 F3 N Ø2
6818- 2C 1H 1	BR F2 SE	7286-1 F2 S	7122- 7 NA 02	6275-1 02 SI	6933-2 N2 O
6568-16	EL F N O	7225-3 N2 6582-7 NA 0	4898- 3C 3H 4 8A N 04	P 7895-1 PB	7437- 3 N2 0
7462- 3C 1H 2	CL F2 O P	4626-22 N4 5	5366-4 CL N3 S	6973- 4 B CL N	7824- 6 N2 0 4864-25 N2 0
5464- 1	03 F SI	4706-1 N4 5	6914-23 CL2 F4 5	I 6973-14 B I N	4864-23 NZ 02
6172-13C 1H 3	BR HG	5912-1 0	6923- 6 CL2 02 5	5449-1 N2 03 P 7137-90 3H12 CLA NA P	S 6523-4 N2 03
6172- 1	CL HG	7575-1 0	4010-34 CL6 HG3	53 6786- 3 P N SI	6680= 1 NA
8018- 1	CLI O P &	1954-1 03 SE	7879- 5 12	7630- 2 FE N 02	52 6523- 8 N4 03
7039- 6	CL2 P SE	5465- 2 BZ D5	5147-7 N NA 02 5147-5 N NA 03	5569-11C 3H13 CL2 N3 Z	R 6523-9 N4 03
5464- 6	D F2 S1	6997-1 BA 04 P	4804- 1 NZ O	6695- 2C 3H20 B10	6465-1 0
5479- 1	F3 S SI	6172-14 BR HG	6937-2 N2 O	7043- 4 B10	4847-1 02
4683-23	13 SM	7569- 2 CL2 02 P	7305- 4 N2 53	6695- 4 CO N7 08	4779-7 53
4682- 3	13 SN	7039- 7 CL2 P SE	6959- 7 N4 03 52	4850-29C 3H22 02 513	4898- 1C - 5 BA F N 04 P
7252- 4C 1H 4	B F6 N	4683-24 13 SN	7669-2 N4 04 S	6524- 1C 3H40 A2 018	4898-2 BAFN04P
6355- 2	CL2 GE	6997- 2 NA2 04 P	7868- 6C 3H 5 8 F2	6524- 3 018 XE2	6364-7 BR N4
6863- Z 7568- 1	CL2 GE	6997- 3 NA3 07 P2	6818- 4 BR F2 SE	7871- 3C 4 B CL F6	4680-1 BR O
5748- 7	F 03 P	7076- 3 CL N S	5045-5 BR2 N 03	6340= 6 CL F6 P	6562-3 CL
6863- 1C 1H 5	BR GE	5529-1 GL 02 P	6292-1 CL HG	7599- 4 CL4 0 52	7758- 2 CL F2 02
5464- 2	FSI	6016- 2 CL 02 P S	6111-15 CL 02 6923-5 CL 05 ST	7599-2 CL4 S2	7090- 4 CL N2
5524- AC 1H 6	B CLA N	7124- 3 CL2 CU N2 012	5826-1 CL2 N 0	7599-1 CL8 52	5278- 2 CL O
6355- 4 6973- 70 1H 7	GE	7568-2 CL2 N O P	7076- 1 CL2 N O	\$ 7599-3 CL8 52	5232- 4 CL 02
6973- 1	BCLN	5748-5 F 03 P	6323-12 CL3 02 3	5472-5 CO HG NA	S2 6033-1 CL 02 S
6973-11	BIN	5748- 2 F2 05 P2	5147-1 N 03	6864- 5 D7 I	7614-18 CL4 N O
6695- 1C 1H15 7043- 1C 1H16	CO N6 06	4682-2 IZ SN 5479-2 MN	7318-1 N 03	6864- 6 D8	4980- 5 CL5 5
7063- 2	B10	6853- 5 O SE	7505-1 06 P	7480-1 F6	5022-1 F3 04
7064- 2	810	4838-9 0 SI	4898- 4C 3H 6 BA N 05	P 7543- 2 F7 N 03	5352- 9 F3 04
6524-10C 1H34	A2 CL4 017	5095-2 BR 02 SI	5956-1 BR CL O 6871-1 CL F3 SH	6340-8 F7 P	7830-10 F3 04
6524-11	CL4 KR2 017	6355- 3 CL GE	5826- 3 CL N O	6165- 4 F9 N O	5216-1 N
6524-12 6524- 7C 1H25	CL4 017 XE2	5095-1 CL 02 SI	7107-7 CL N O	6165-7 F9 N O	7255-18 N O
6524- 8	CL3 KR2 017	7122-17 N	7044- 2 CL2 N R	52 5023- 4 F10 N2	6111-14 N O2
6524- 9	CL3 017 XE2	5170- 3 N G	6177-1 CL2 SI	7246-1 F12 HG (	D2 P2 5216- 5 N S
6524- 4C 1H36	A2 CL2 017	4658-27 02 P S2 4858-2 03 P S	6914-22 CL2 SI	5023- 3 F12 N2	6364-1 N3
6524- 6	CL2 017 XE2	5196- 6 04 P S	4625-1 N NA 04	5 6474-16 K3 N3 03 5 6474-12 K3 N3 04	5 6523-5 N3 02
6340- 4C 2	AS CL2 F3	5524- 3C 2H 8 8 CL4 N	7612- 1 N NA 05	5 86474- 8 N3 NA3 (	3 \$ 7669-1 N3 02 5
7871- 1	8 CL2 F3	7761-3 BN	7581-3 N2 52 4706-4 N4 5	6474- 4 N3 NA3 (	04 6523-10 N3 03
7252- 2	82 8A F12	6493- 2 BE FZ N2 04	5859- 2 N4 5	7199- 4 CL F2 N	2 7669-4 N3 05 S
5465- 3	82 010	6031- 3 BR2 N4 S2 TE	7305-1 44 52	7199- 5 CL2 F N	6523- 2 N5 02
6493- 1	BE F2 K2 04	6031- 2 CL2 N4 S2 TE	5954- 2 03 SE	7199-6 CL2 F N	7868- 3C 4H 6 B CL
6493- 3	BE F2 NA2	7224- 1 CL6 N6 P4	6923-8 06 52	6474-15 K2 N3 0	3 5 5352-16 BR F3 02
6853- 2	BR CL FA SE	7500- 4 CO N6 S2	6923- 9 07 S2	6474-11 K2 N3 04	5812- 3 BRZ 03
7076-18	BR2 CL2 N2 .52	6355- 5 GE	6328- 4C 3H 7 8 BR2	6474- 3 N3 NA2	73 5 5480- 2 BR2 03 5
7303-27	CA F2 05 5	7500- 6 MN N6 52	7868-5 B F2	6474-28C 4H 2 AG N3 04	5052-2 CL N5
7303- 8	CL F3 03 S	4756-1 N D3 P S	4961-1 BA N2 0	7742-1 BCL3 F	02 5243-1 CL3 04 P
7303- 3	CL2 F2 03 5	7500- 3 N6 NI 52	6172- 3 CL HG	7742- 3 BR CL2 1	13 5243- 6 CLS 04 P
7303- 4	CL2 F2 03 5	7500- Z N6 S2 ZN	6627- 2 CL HG O	7501- 5 CL N 03	6864- 3 D I
7303- 6	CL2 FZ 03 S	6973- 9C 2H 9 B BR N	6787-4 CL 02 S	7303-36 CL N3 0	6864-4 D2
7303-19	CL2 F2 06 52	6973- 3 B CL N	7569- 3 CL2 02 1	4839- 3 CL5 F3	51 5352-5 F3 N 03
6853- 4	CL2 F4 52	7730-23C 2H10 N4	5845-2 K 04 S	7597-6 CL8 52	5352- B F3 N 03
7076-20	CL2 N2 S	6007-11 0 SIZ	5826- 2 N O2	7199- 2 F2 N2	5147- 6C 4H 6 N NA 03
7199-1	CL4 F2	6122- 1C 2H14 B2 N2	6484-3 N S	7199-3 F2 N2	4804- 2 NZ O
7814- 2	CL5 F3 S1	7043-3 B10	5326-1 06 P	6474-14 K N3 03	6697+1 N2 O \$ 7025-20 N2 07
5023- 7	F3 N	7063-1 B10	5529- 2C 3H 8 CL 02 P	6474-10 K N3 04	7255-16 N2 02
7543- 1	F3 N 03	7063-3 810	7568- 3 CL2 N 0	P 7753-1 K N3 05	5283-10 NZ 0Z 52
5023- 8	F4 O	7064-1 810	6914-21 CL2 SI	6474- 6 N3 NA 0	3 5 6364- 8 N4
7303- 1	F4 03 5	6695- 3 CO N7 OB	7225- 2 N SI	6474- 2 N3 NA 04	7730-18 N4
7303-14	F4 03 5	4821- 6C 3 BR CL2 F2 4819- 3 BR2 K N3 O3	4802 1 N4 02	4901- 2 02 S3 5352- 10 AH 3 BP F3 H	6523-1 N4 0
7303-15	F4 06 52	7199- 8 CL F2 N3	4705-1 N6 53	7437- 4 BR N2 0	6523-11 N4 02
6340- 7	F5 P	6165-1 CL F6 N O	4838- 8 0 51	4631-1 BR N2 S	6523-12 N4 02
6853- 3	F6 SE2	7199-9 CL2 F N3	6923-1 06 52 6923-2 07 52	6364-2 BR2 N3	2 6959-8 N4 03 52
5023- 2	F7 N	6165- 2 CL2 F5 N O	4683- 2C 3H 9 BR SN	6711-17 CL N2	7669- 5 N4 04 S
7814- 4C 2H 1	CL4 F3 51	4819-2 CL2 F5 N 0	7250-1 BR3 N P	7051-1 CL N2 0	6967- 2 N6 02
6818- 3	F2 N SE	4819- 1 CL2 N3 03	4850-28 CL N 0	6940- 2 CL N4 0	2 5803-1 NE 52
7303- 9	F3 03 5	4821- 5 CL3 F2	4683- 1' CL SN	7079- 1 CL N4 0	6712-1 0
7303-24	F3 04 5	7303-41 F3 N 03 52	6166- 2 CL3 N3	3 5079-83 CL 02	4668-10 02
7303-11	F3 06 52	7199- 7 F3 N3	7869-1 F SN	4839-1 CL4 F3	SI 5480-1 03 S
7303-21	F5 02 5	7303-12 F6 03 S	7630- 1 FE 02 5	6864- 8 D5	6033- 6 03 5
6959-13	CL N3 S	7303-17 F6 06 52	4683-3 I SN	4999-1 N3 04	5910-2 03 52 4010-18 53
7814- 3	CL2 F2	7303-18 F6 06 52	5479- 3 LI MN	6523- 7 N3 04	
7814- 5	CL3 F3 SI	7856-2 F7 N	5170- 4 N O	6940-1 N3 04 7310-7 N5 F	
6914-26	CL4 GE	6165- 9 F7 N G		·····	
6914-27	CLA SN	6165-10 F7 N D			
6818- 1	F4 SE2	5023- 5C 3H1 F6 N 0			
6737- 1	N2 0		Fig. 1	Old format of annulation	o molecular (
3910- 1	05		rig. 1.—		e molecular formula index

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Fig. 2.—New format of cumulative molecular formula index

P	571 573		P			P			D		
(P2 (L3)	1274 1P2 CO23	1275	CONTINUES 1276 (PZ N)	[P2 N2]	1278 (P2 NA3)	CONTINUED 1279 (P2 05)	1280	1281 (#3 GD3)	CONTINUED 1282	1283 P3 N9	1284 (P4 CL)
N 02 PT2 S	#7 CO3	P2 811	NI 5 C23 H37	08 IN C36 H30 010 C8 H18	07 C2 H5	S2 C12 H28 S12 C10 H28	P2 015	H6 N 012 H12 N3 012	P3 F12	C6 H24	OS C12 H33 OS C20 H49
N2 02 C12 H11 N2 02 C14 H15		P2 H14	0 C30 H25	010 C24 H34	P2 845	P2 06	V3 C70 H90	H15 N4 012	C7 H9	C12 H24 C12 H36	05 C58 H55
N2 02 C16 H19 N2 03 C19 H15	PZ CA	N4 06	02 PT C7 H21 02 PT C13 H33	010 C32 H50	P7 NI	C4 H12	P2 025	P3 CL	P3 FE	C18 H48 C24 H60	RU C12 H33 RU C20 H49
0 05 C37 H30 0 PT2 5 C25	H10 N2 06	P2 H17	03 PT C7 H21 03 PT C19 H37	H36	C16 H32	C6 H16 C7 H16	C44 H92	IR C54 H47	012 C12 H30	C36 H36	RU CSO H45
0 RE C10 H24	04 C14 H24 04 C18 H24	N5 06	03 PT C21 H41 04 C9 H23	H38 010 52 C18	C18 H36 C24 H40	C7 H18 C8 H16	P2 049	N3 C10 H25 N3 C15 H35	012 CZ4 H54	C36 H84 C42 H48	P4 CL2
0 RE C12 H30	04 C24 H30 D4 C29 H22	P2 H36	04 C16 H21 05 C7 H19	H34 010 52 C22	C28 H40 C30 H52	C8 H20 C9 H22	C92 H148	N3 C20 H45 N3 C30 H25	P3 FE2	06 C24 H48 06 C30 H36	CU 012 C12
0 RE C34 H30	04 C40 H30	N12 N13 SE8	05 C10 H23 06 C5 H15	010 S2 C39	C34 H52 C36 H40	C10 H22 C10 H24	P2 P0	N3 C40 H85 N8 C10 H20	015 C45 H33	06 56 C18 H36 015 C57 H48	CU 016 C12
02 RE C36 H30	G10 C40 H30	P2 HG	04 C8 H19	010 S2 C40	C38 H44 C38 H52	C11 H18 C11 H24	C14 H36 C24 H40	NB C30 H30 NB C35 H40	P3 H3	D18 C36 H24 56 C6	FE C12 H32 FE C20 H48
PT2 5 C24 H53 RE C36 H30	P2 CR2	14 C8 H24 04 54 C4 H12	06 C9 H21	010 52 642	C44 H48	C12 H20	C28 H30	0 05 C31 H46	NG	P3 N12	FE C28 H48 NJ 012 C72
RU C36 H30	N12 SE SN C48 H52	04 54 C8 H20 04 54 C12 H28	06 C10 H23 06 C10 H25	010 S2 C46	NZ C42 H36	C12 H28	C38 H36	0 40 C55 H46	F3 H12	021 CS7 H45	05 C12 H32
P2 CLA	P2 654	06 S2 C8 H20	06 C12 H27 06 5 C12 H27	010 U C24 H54 011 C9 H14	02 C12 H24 02 C16 H24	C13 H28 C13 H30	PZ PD2	51 C6 H18	P3 13	P3 N15	05 C28 H48
NR C38 H36 NZ 02 C4 H8	H024 083 112	P2 H04	07 C5 H13 07 C7 H17	012 C9 H14 012 C14 H25	02 C24 H23 02 C28 H24	C14 H30 C14 H32	SE C30 HT0	P3 CL2	09 RE C34 H45	56 C6 H18 56 C30 H66	05 652 HAB
N2 08 C23 H28	P2 CU	010 EN3	07 C9 H21 07 C11 H25	012 U C24 H54 013 5 C28 H52	02 C32 H24 02 C38 H30	C15 H32 C16 H20	P2 P1	F4 N3 IR C54 H46	09 RE C63 H63	54 C42 H42	RU C12 H32 RU C20 H48
N2 52 C4 H8	1 04 C24 H38	P2 1	07 C12 H27 07 C13 H29	014 U C12 H30 016 C16 H26	02 C42 H38 54 C12 H28	C16 H36 C17 H38	C36 H32	N3 C24 H20 N3 C24 H30	P3 IR	P3 HZL	RU C28 H48 RU C50 H44
HE 04 C2 H12	1 06 C24 H30	C47 H37	07 C15 H17	PD C28 H50	54 C16 H30 54 C16 H36	C18 H24 C18 H40	C36 H48 C38 H36	N7 C8 H16 N7 C8 H24	C54 H48 02 C56 H50	56 C42 H48	RU C52 H48
0 52 02 C1 H2	N 06 C7 H18 N 06 C13 H30	HN N2 D8 C36	07 C19 H25	FZ N3	54 C32 H48	S C10 H24	P2 P12	N7 C12 H32 N7 C16 H40	04 C56 H48 04 C57 H50	P3 NI	P4 CL3
02 C2 N4 02 S	N 06 C19 H42 N 06 C25 H54	N 04 C10 H26 NI C30 H37	07 C23 H33 07 C25 H21	0 RE 53 C23	#2 0	S C16 H36	54 C26 H62	N7 C28 H32	06 L38 H32	0 C10 H27	RE 652 H48
03 C4 H8 04 S2 SH C6	N2 04 C30 H34 N2 08 C34 H30	PT C7 H21 PT C24 H39	07 C27 H25 07 \$ C14 H31	0 RE 53 C39 H30	C31 H24	5 C20 H44	P2 RE	P) (1)	N3 021 C34	010 C19 H45	NA (24 H20
H18 RE C36 H30	P2 0	P2 12	08 C34 H33 015 C40 H47	02 C2 H7 02 C3 H9	P2 02	52 ZN C8 H20 54 C14 H20	C36 H33	F3 N3	нат	010 C37 He1	N8 C20 H40
SE7 SN C34 H30	N 06 C8 H20	C12 H10	017 C14 H25 017 C14 H27	04 C15 H25 05 S C38 H31	C5 H14	54 C15 H22 54 C16 H24	PZ SZ	N3 C3 H9 N3 C18 H15	P3 L14	P3 0	Po CLA
P2 CL5	P2 02	C17 H24 C16 H26	PD C14 H33 PD 5 C14 H33	07 C5 H21 011 C9 H15	C6 H16 C10 H24	54 C17 H26 54 C20 H32	C4 H12 C6 H16	N3 03 C18 H15 N3 03 C30 H21	N5 010 C7 H8	C58 H31	FE2 C20 H48
1 NI C18 H30	07 C15 H26	C20 H30 C28 H30	PT C10 H25	012 C6 H13	C13 H28 C17 H38	TH C14 H14	CB H20 C10 H24	N6 C6 H12 N6 C12 H30	P3 H0	P3 03	FE2 C28 H48 N6 C2 H8
P2 CLA	P2 07	1R 0 C37 H30	\$ C30 H25	PZ -N4	C18 H40 C19 H42	P2 07	C12 H28 C14 H16	N6 C18 H42 N6 C25 H54	03 C12 H27 03 C21 H45	¥ C44 H39	N6 C4 H12 N6 C8 H20
FE2 04 PT C50	07 C5 H5	NN 02 C34 H30	PZ NZ	02 CB H24	C22 H26	C3 H10	C16 H2C	N6 03 C45 H36	03 C29 H33 03 C37 H33	P3 04	N6 C10 H20 N6 C12 H12
12 NI C6 H12	P2 F	N2 NI2 02 C36 H30	C38 H28 C42 H34	04 54 C14 H16	C25 H22	C5 H12	C19 H26	N9 06 C66 H60	03 C57 H45	C18 H15	N6 02 C8 H16
12 RE C36 H24 N2 C2 H6	06 C6 H15 010 C6 H13	N2 PD C30 H56 NI C6 H18	NI C24 H21 NI D2 C36 H30	07 52 C10 H28 08 66 H16	C26 H24 C27 H26	C6 H16	C21 H30 C22 H32	RE C30 H45	012 C21 H45	502 C18 M18	03 32
N2 HE SZ C38 H24	P2 F2	NE C14 H24 NI C28 H46	NI 06 C6 H18 NI 06 C28 H46	08 C26 H24 D10 C14 H25	C28 H60 C31 H66	C8 H20 C10 H16	C24 H20	P3 CL4	P3 N	P3 010	NA CLU
N2 02 C8 H4 N2 03 5 C7 H4	H2 05	NI C32 H38 NI C42 H42	NI 06 C32 H38 NI 06 S2 C44	.010 C16 H20 015 C20 H28	PT C26 H44 RE C40 H45	C10 H20 C10 H24	C28 H28 C28 H40	F2 N3 HG2 03 C54	0 RH C34 H45	C12 H13	P4 (1.1.
NZ PT C42 H34	N 05 C10 H25 N2 03 C8 H20	NI 02 C36 H30 NI 06 C42 H42	NI 08 C36 H30	015 C39 H42 017 C20 H30	5 C8 H20 -5 C12 H28	C12 H10 C12 H12	C29 H42 C30 H44	H39 N5 C2 H8	014 U C40 H90	P3 012	N3
NI C18 H30	NZ 05 C14 HZ0	05 C6 H14	NI 52 C8 H18 NI 52 C14 H30	018 RU C24	S C16 H20 S15 C40 H44	C12 H28 C14 H16	PZ 54	N5 C4 H8 N5 C4 H12	P3 N2	PR C24 R34	P4 0.14
02 C1 03 C14 H10	N2 05 C20 H32	RE C42 H42	NI 52 C30 H46	PT2 54 C22 H42	P2 03	C14 H28 C14 H32	C12 H28	N5 C8 H20 N5 C16 H36	011 C18 H39 014 C9 H15	P3 014	HO6 04 C72
PT C36 H34	H18	PZ (3	NI 52 C44 H42	S2 C9 H24	C14 H16	C16 H20	C15 H30 C16 H20	N5 C24 H44	015 C9 H15	C6 HL7	H62
P2 CLU	NA2 05	IR 04 C36 H30	02 C24 H18	PZ N5	B2 04	C18 H40	C18 H32 C17 H34	P3 (L)	P3 N3	P3 PT	P4 C0
C6 H5	05 C2 H6 05 C4 H10	P2 14	02 5 C36 H30 02 55 C18 H16	09 C10 H13	C4 H12	C20 H44	P2 S1	N4 C2 H6	C18 H42	C54 H47	C52 H48
C7 H7 C8 H7	05 C8 H18	N RE C44 H37 RE C36 H30	03 C24 H18 03 PT C24 H42	010 C10 H15 010 C14 H23	C6 H16 C8 H2G	C28 H28 C28 H60	C10 H26	P3 CLA	C36 H30	P4	P4 C02
03 V	P2 F3	P2 K	03 S2 C20 H22 04 C8 H22	010 C16 H19 010 C17 H21	C10 H16 C12 H10	C30 H25 C32 H68	P2 515	N3	ND 021 C36	C8 H20 C24 H44	H34 N10 012
P2 CL10	C4 H9 C13 H27	D6 C16 H17	04 C14 H30 04 C14 H34	010 C22 H23 010 5 C12 H19	C14 H16 C16 H20	C36 H76 C40 H84	C42 H48	N3 C8 H8 N3 U6 C36 H24	06 C6 H18 06 C12 H30	P4 AG	P4 CR
FE C32 H30	#2 #4	P2 L12	04 P0 C28 H38	011 C28 H35	C16 H28 C18 H36	SIZ CB HZ4 SIZ C10 H28	P3	N6 09 C87 H45 N9 C30 H30	04 C18 H42 06 C24 H54	CL 04 C72 H60	02 C30 H48
02 C4 H4	C2 H2	C12 H10	04 S8 C15 H36	012 C31 H45	C24 H28	72 C15 H26 22 D8	C26 H33	N9 C36 H90 N9 C54 H126	06 C36 H30 06 C42 H42	N 03 C72 H60	02 C52 H44 02 C54 H48
02 ¥	842 NI C6 H12 CL2 NI C6 H12	C15 H28 C16 H18	04 S8 C22 H36	015 C16 H23	PB 54 C4 H12	C20 H24	C41 H39	H42	021 Y8 C36	P4 84	P4 F8
P2 CL11	12 NI C6 H12 N2 NI D6 C6	C16 H30 C17 H20	04 58 C42 H34 06 C4 H14	016 C16 H25	52 C4 H12 52 C6 H16	C48 H42 RE2 C44 H30	P3 AG	H39	025 (23 H38	CLIS NO CO	N4
C2 H3 C3 H5	N2 NI 52 C8	C17 H32 C18 H22	06 C12 H12 06 C12 H14	P2 N6	52 C8 H20 52 C12 H28	52 C20 H40 52 C21 H42	8 F4 09 53 C9	P3 619	56 C12 H30	FA 543	N1 C24 H20
C4 H7	H12	C18 H34 N2 017 C15	06 C12 H26 06 C12 H30	NI 010 V2 C82 H54	52 C15 H34 52 C16 H36	52 C22 H44 513 C9 H18	09 53 C12 H32	N& 09 C57 H42	56 C24 H54 56 C36 H30	N4 022 C18	P4 F12
	P2 F12	N3 017 C17	06 C18 H18 06 C23 H30	0 C10 H30 0 C30 H30	53 C4 H12 53 C8 H20	28 CI4 H18	P3 AL	P3 CL12	56 C42 H42 56 C48 H102	N10 020 52	NI
HUG U2 C36 H30	0 64	N5 016 C16	06 C24 H20 D6 C24 H22	02 C12 H24 07 C30 H26	53 C12 H28 54 C2 H8	P2 09	09 C24 H30	N	56 C72 H150 56 C96 H198	C24 H32 D14 C10 H18	P4 FE
P2 CD	P2 F18	#2 1 T 3	06 PB C30 H34	09 C14 H20	54 C6 H12	S12 C4 H12	P3 82	P3 CL13	P3 N4	P4 84	C28 H50 I C20 H49
C30 H52	C12	NZ 014 C15	D6 52 C18 H36	S2 C12 H24	54 C8 H20 54 C9 H22	#2 010	C42 H77	03 11	018 C20 H29	RU C12 H33	16 C72 H60
C42 H48 C44 H48	P2 FE	07 C5 H9	06 52 C30 H52	P2 N7	54 ZN C24 H52	C6 H14	F3 63		UI4 C58 M57	RU C20 H49	P4 H2
CR N6 03 54 C19 H36	12 C36 H30 H3 09 C36 H30	P2 MN	07 C10 H12 07 C14 H20	012 C19 H23 013 5 C21 H35	W C14 H24 W C16 H30	SIZ C12 H26	C10 H32	P3 (02	C10 H30	F9 842	NZ 3
F12 C11 H11 12 C36 H30	N3 53 C39 H30 D3 C39 H30	NZ 08 C36 H30	07 C14 H34 07 C18 H28	015 C19 H25 015 C23 H31	W C18 H24 W C24 H30	U2	C12 H36 C24 H60	H27 NB 012	010 C9 H16	NI C28 H48	29 H3
12 02 C36 H30 H2 02 52 C38	PZ H	P2 MH2	07 C22 H36 07 C25 H40	015 C23 H33 016 C24 H27	W C29 H22 W C30 H24	P2 011	C117 H102 BR6 C12 H30	P3 CA	013 C10 H16	RU C20 H48	PA HA
H30 N2 08 C36 H30	N NA4 06	N6 NI 08 C80	08 C10 H12 08 C12 H26	P2 NB	P2 05	C16 H36	CL6 C6 H18 CL6 C12 H30	H19 N4 D9	013 CT2 H20 014 C29 H42	P4 CA	NR
N2 52 538 H30	N 06 TL4	D8 C22 H24	08 C14 H20 08 C16 H20	02 C20 H28		P2 012	16 C12 H30	03 C44 H39 012 C39 H81	015 C14 H24	H14 018 07	P4 H6
P2 C02	010 SHA	P2 M0	08 C18 H28	02 5 C12 H24 04 52 C18 H20	C2 H8	C6 H14 C12 H22	P3 BAZ	012 CS7 H45 015 S3 C27	P3 N6	P4 CD	N2 53
CD 06 C42 H30	#7 H3	04 C14 HZ4	08 C20 H28	022 L96 H46	C8 H20	C42 H46	NZ 014 C9 H11	H48	C9 H27 09 C57 H51	FE 14 C72 H60	NZ SE3
HG 06 C42 H30	K2 N 06	04 C18 H24 04 C24 H30	08 C20 H48	P2 N11	C12 H12	PB3 C28 H16	F3 0K	P3 C#3	U22 C30 H39	14 MN C72 H60	P4 H9
H54 N6 H1 012 C76	P2 H4	04 C29 H22 04 C30 H24	08 C23 H32 08 C24 H28	013 624 H32	C12 H28	2N C44 H52	0 05 C55 H46	P3 03	P3 N7	P4 C05	N3 53 N3 563
04 C54 H40	N2 06	010 C10 H18 010 C16 H30	08 C24 H36 08 C24 H50	P2 NA	C17 H28 5 C12 H24	P2 013	P3 843	IR C54 H44	017 C21 H28	D4 C1	P4 H12
06 C18 H30 D6 C42 H30	06 52	010 C40 H30	08 C27 H40 08 C28 H44	07 C3 H9	S C13 H28 S C15 H34	C81 H166	05 C54 H45	P3 F6	P3: NE	CU 04 C22 H/O	N4 53
D4 C42 H66 D6 P8 C42 H30	P2 H6	P2 H	D8 C29 H40 D8 C31 H48	P2 NA3	52 C6 H16 52 C8 H20	P2 014	P3 C03	N3	017 5 C31 H43 026 C59 H67	FE C12 H33	P4 H13
D6 ZN C42 H30	NZ 06	NA3 04 C6 H6	D8 C32 H52	27 2012	52 C10 H24	C10 H18		N3 C24 HZ4		FE C28 H49	

Fig. 3.—Rotaform Index

ore curbe	on aloni	s and s of more oxyden alon	-				
I.C. 859	L.J. Ame J. P	COID BORATES I. STRUC Leeson, J.A. Lowery, G rican Cyanamid Co., Pea harm. Sci. 50(3), 193-7(19	TURAL CON G.M. Sieger, rl River, N. 961).	ISIDERA , S. Mul Y. Recd	[10NS. 1er. . Mar. 16, 1	1960.	
		снон					
			lla—Carbons 1,2 11b—Carbons 1,3	: CHCH. CHCH			H H
	1)	C21 H28 B F O7	(lla)	(as l	Na sal t)	Fr U	1
	2)	C42 H55 B F2 O12	(111a)		**	We Color 18. CH - CH (F 8 Be	<b></b>
	3)	C21 H26 B F O7	(IIb)	(as l	NH <sub>A</sub> salt)	IIIb-Carbons 1,2: CH-CH (F-4 Bon	ites)
	4)	C42 H51 B F2 O12	(ШЬ)				
Search	2:	Sulfanilamidodiazines (dia	zine = 6-mem	bered rin	g containing t	wo nitrogen atoms).	
a. Wit	th or w	ithout alkyl substituents. I	Jnder SN40	2 search	C10H10. Fo	r alkyl substituents search	
<sup>C</sup> 10 + I	n <sup>H</sup> 10 +	2n*					
			1				
	20)		R=				
	20)	C12 H14 N4 O2 S	F+				
	21)	C12 H14 N4 O2 S	Dr.		/=	- NH-SO2- NH2	
	22)	C14 H18 N4 O2 S	:_B.		<u> </u>		
	23)	C14 H10 N4 O2 S	n-amyl		Π-	-*	
	24)	4245	in any i		m.f.		
L 16.	I.C.	4343	der SNLO	earch C	н.с.	10010-3	
<b>D</b> . IF	emer si		403	searchic	11 + n ''12 +	- 2n*	
	N		T				
	1)		Y=			NH-SO2-	
	2)	C12 H14 N4 O3 S	F+O			Lon	
	2)	C13 H16 N4 O3 S	PrO			N <sup>-</sup>	
	4)	C14 H18 N4 O3 S	BuO		5)	C11 H12 N4 O3 S	R= Me
	5)	"	sec-BuO		3) 7)	C12 H14 N4 O3 S	Et
	1.C.	10035			1.C.	6360	
c. Fa	or dieth	ners, under SN101, searcl	hC. H.				
		4 4 R1	12 + n	14 + 2n°			
			$\begin{bmatrix} R_2 \\ R_3 \end{bmatrix}$				
			R <sub>l</sub> =	<sup>R</sup> 2 <sup>=</sup>	R <sub>3</sub> =		
	1)	C12 H14 N4 O4 S	MeO	MeO	н		
	2)	C13 H16 N4 O4 S	EtO	и	н		
	12)	н	MeO	Me	MeO		
	13)	C15 H20 N4 O4 S	EtO	11	EtO		

Search 1: Borate derivatives of steriods. Under B headings examine all compounds containing 19 or more carbon atoms and 3 or more oxygen atoms.

I.C. 12534

Fig. 4.—Two searches using the RotaForm Index

The user must in each case decide which is less costly additional filing space and card processing (repeating structural diagram under each element) or additional clerical time in looking up diagrams from a serial number. In the former case one also has to consider cost of refiling cards if they are to be shown to chemist or cost of reproducing them. In the latter case it is the time to write down serial numbers.

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