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Intermediates in the 1,2-(NH₂)₂C₆H₄/P(NEt₂)₃ Transamination Reaction.

Page 4313. The parameters of C(20) and C(20') are incorrect in Table III. The correct parameters are as follows:

C(20)	-0.0841 (14)	1.1865 (24)	0.2144 (9)	0.144 (10)
C(20')	-0.1666 (13)	0.1929 (22)	0.1045 (14)	0.165 (12)

Page 4314. All parameters in Table IV are incorrectly multiplied by 1.414. The corrected table is given below.

Table IV. Atomic Coordinates^a ($\times 10^4$) and Equivalent Isotropic Displacement Parameters ($\text{\AA}^2 \times 10^3$) for [C₆H₄(NH)₂P(S)]C₆H₄(NH)NP(S)NEt₂

	<i>x/a</i>	<i>y/b</i>	<i>z/c</i>	<i>U</i> (eq) ^b
P(1)	-383 (1)	-571 (1)	8738 (1)	35 (1)
P(2)	-9 (1)	1945 (1)	8102 (1)	37 (1)
S(1)	630 (1)	-607 (1)	9308 (1)	47 (1)
S(2)	733 (1)	1110 (1)	7943 (1)	49 (1)
N(1)	-790 (2)	-2005 (3)	8294 (2)	45 (2)
N(2)	-890 (2)	-532 (3)	9114 (2)	41 (1)
N(3)	-602 (1)	806 (3)	8186 (1)	35 (1)
N(4)	-656 (2)	2684 (4)	7420 (2)	52 (2)
N(5)	300 (2)	3070 (3)	8727 (1)	43 (1)
C(1)	-1303 (2)	-2554 (4)	8453 (2)	41 (2)
C(2)	-1357 (2)	-1703 (4)	8937 (2)	40 (2)
C(3)	-1838 (2)	-2009 (4)	9170 (2)	50 (2)
C(4)	-2253 (2)	-3224 (5)	8914 (2)	65 (2)
C(5)	-2200 (2)	-4076 (5)	8436 (2)	65 (2)
C(6)	-1723 (2)	-3754 (4)	8191 (2)	53 (2)
C(7)	-1324 (2)	1150 (4)	7694 (2)	40 (2)
C(8)	-1332 (2)	2233 (4)	7266 (2)	46 (2)
C(9)	-1966 (2)	2753 (5)	6748 (2)	65 (2)
C(10)	-2589 (2)	2152 (6)	6676 (2)	75 (2)
C(11)	-2581 (2)	1091 (5)	7101 (2)	68 (2)
C(12)	-1951 (2)	570 (4)	7613 (2)	51 (2)
C(13)	-188 (2)	3903 (4)	8880 (2)	60 (2)
C(14)	1058 (2)	3428 (4)	9130 (2)	53 (2)
C(15)	-253 (3)	3376 (5)	9487 (3)	87 (3)
C(16)	1281 (2)	4647 (5)	8823 (2)	82 (3)
P(1')	5496 (1)	-294 (1)	4122 (1)	33 (1)
P(2')	5058 (1)	2307 (1)	3190 (1)	37 (1)
S(1')	4490 (1)	-440 (1)	3789 (1)	50 (1)
S(2')	4280 (1)	1527 (1)	2372 (1)	53 (1)
N(1')	5936 (2)	-1628 (3)	4006 (2)	41 (1)
N(2')	6004 (2)	-354 (3)	4960 (2)	39 (1)
N(3')	5681 (1)	1174 (3)	3794 (1)	34 (1)
N(4')	5680 (2)	3059 (4)	3051 (2)	48 (2)
N(5')	4788 (1)	3394 (3)	3587 (1)	42 (1)
C(1')	6409 (2)	-2305 (4)	4612 (2)	37 (2)
C(2')	6446 (2)	-1570 (4)	5166 (2)	37 (2)
C(3')	6883 (2)	-2043 (4)	5817 (2)	52 (2)
C(4')	7274 (2)	-3274 (5)	5897 (2)	68 (2)
C(5')	7238 (2)	-4006 (5)	5345 (2)	65 (2)
C(6')	6801 (2)	-3531 (4)	4688 (2)	49 (2)
C(7')	6393 (2)	1624 (4)	3960 (2)	38 (2)
C(8')	6370 (2)	2712 (4)	3520 (2)	43 (2)
C(9')	6986 (2)	3333 (4)	3585 (2)	56 (2)
C(10')	7626 (2)	2864 (5)	4100 (2)	65 (2)
C(11')	7650 (2)	1813 (5)	4544 (2)	65 (2)
C(12')	7033 (2)	1180 (4)	4481 (2)	51 (2)
C(13')	4058 (2)	3868 (5)	3348 (2)	61 (2)
C(14')	5317 (2)	4115 (5)	4209 (2)	64 (2)
C(15')	3903 (3)	5295 (5)	2994 (3)	112 (3)
C(16')	5385 (3)	3460 (5)	4856 (2)	108 (3)

^aAtoms have occupancies of 1.0. Atoms in molecule B are primed.

^bThe equivalent isotropic *U* is defined as one-third of the trace of the orthogonalized *U_{ij}* tensor.

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