

**Supplementary Table S3. Phospholipids identified in the “oxysterol window” of the plasma metabolome**

Measured <i>m/z</i> (RA <sup>a</sup> )	Retention time/min	Elemental composition	Lipid Maps hit <sup>b</sup>	Metlin Library hit <sup>c,d</sup>	Measured MS <sup>2</sup> fragments (RA)	Postulated structure
494.3240 (3.75)	9.52	C <sub>24</sub> H <sub>49</sub> O <sub>7</sub> NP	GPCho 16:1(9Z)/0:0	GPCho 16:1(9E)/0:0 <sup>d</sup>	476 (100), 417 (0.86), 184 (33.4)	GPCho 16:1/0:0
496.3402 (100)	10.57	C <sub>24</sub> H <sub>51</sub> O <sub>7</sub> NP	GPCho 16:0/0:0 GPEtn 19:0/0:0	GPCho 16:0/0:0 <sup>d</sup>	478 (100), 419 (0.30), 313 (0.13), 258 (0.52), 184 (29.27)	GPCho 16:0/0:0
520.3400 (58.26)	10.02	C <sub>26</sub> H <sub>51</sub> O <sub>7</sub> NP	GPCho 18:2(9Z,12Z)/0:0	GPCho 18:2(9Z,12Z)/0:0 <sup>d</sup>	502 (100), 443 (0.26), 337 (0.11), 258 (0.61), 184 (29.60)	GPCho 18:2/0:0
522.3553 (32.38)	10.99	C <sub>26</sub> H <sub>53</sub> O <sub>7</sub> NP	GPCho 18:1(9Z)/0:0	GPCho 18:1(9Z)/0:0 <sup>d</sup>	504 (100), 339 (0.47), 258 (0.83), 184 (28.21)	GPCho 18:1/0:0
524.3711	12.52	C <sub>26</sub> H <sub>55</sub> O <sub>7</sub> NP	GPCho 18:0/0:0	GPCho 18:0/0:0 <sup>d</sup>	506 (100), 447	GPCho

(23.39)					(0.37), 341 (0.30), 258 (0.46), 184 (24.53)	18:0/0:0
544.3395 (9.34)	9.95	C <sub>28</sub> H <sub>51</sub> O <sub>7</sub> NP	GPCho 20:4(5Z,8Z,11Z,14Z)/0:0	GPCho 20:4(5Z,8Z,11Z,14Z)/0:0 <sup>d</sup>	526 (100), 467 (0.25), 258 (0.32), 184 (27.49)	GPCho 20:4/0:0
568.3400 (1.26)	9.90	C <sub>30</sub> H <sub>51</sub> O <sub>7</sub> NP	GPCho 22:6(4Z,7Z,10Z,13Z,16Z,19Z)/0:0	GPCho 22:6(4E,7E,10E,13E,16E,19E)/0:0 <sup>c</sup>	550 (100), 385 (3.03), 184 (19.43)	GPCho 22:6/0:0
570.3556 (0.56)	10.34	C <sub>30</sub> H <sub>53</sub> O <sub>7</sub> NP	NA	NA	NA	GPCho 22:5/0:0

<sup>a</sup> Relative abundance as determined by peak area.

<sup>b</sup> From a search of *m/z* value against the Lipid Maps<sup>S4</sup> database <http://www.lipidmaps.org/tools/index.html>

PGCho x:y/x:y is an abbreviation for glycerophosphocholine where the number of carbons in the acyl chains are indicated by x and the number of double bonds by y.

<sup>c</sup> From a search of *m/z* against the METLIN metabolite library<sup>17</sup> <http://metlin.scripps.edu/index.php>.

<sup>d</sup> From a search of M against the METLIN metabolite library<sup>17</sup> <http://metlin.scripps.edu/index.php>