

Supplementary materials

Primary and secondary organic aerosol in summer of Beijing, China

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Fig. S3 Back trajectory clusters during the campaign at (a) CP (b) PKUERS

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Fig. S5 Three categories of PM_{2.5} chemical compositions according to back trajectory clustering analysis

Fig. S6 Primary organic matter from different directions of CP and PKUERS

Fig. S7 Daily average concentrations for different classes of primary organic matter, (a) n-alkanes (b) saccharides (c) PAHs (d) hopanes

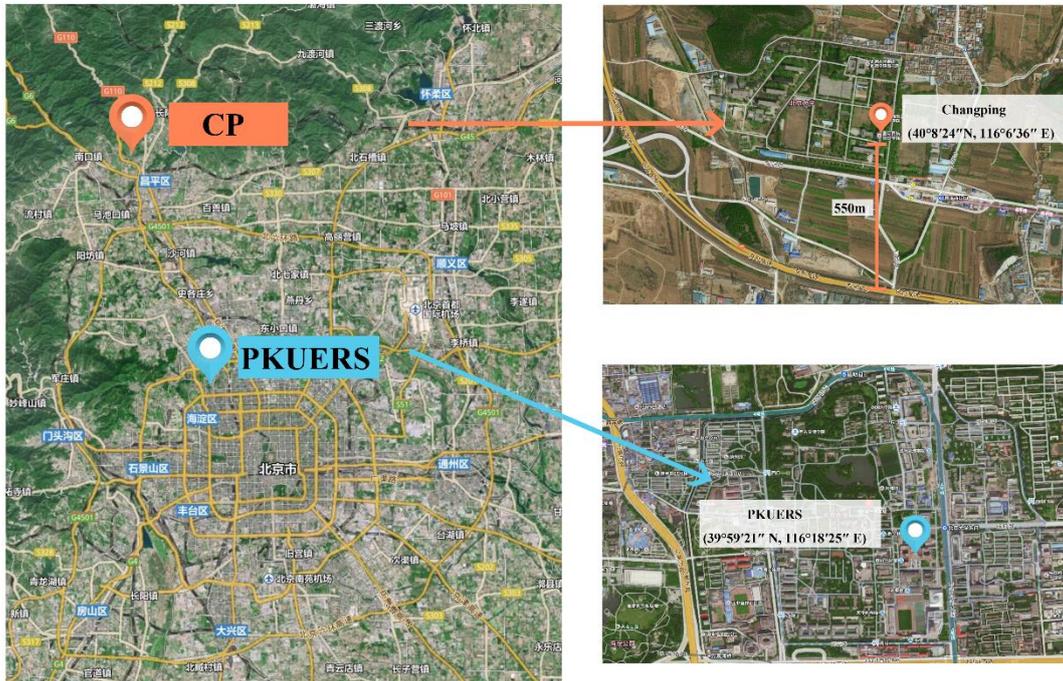


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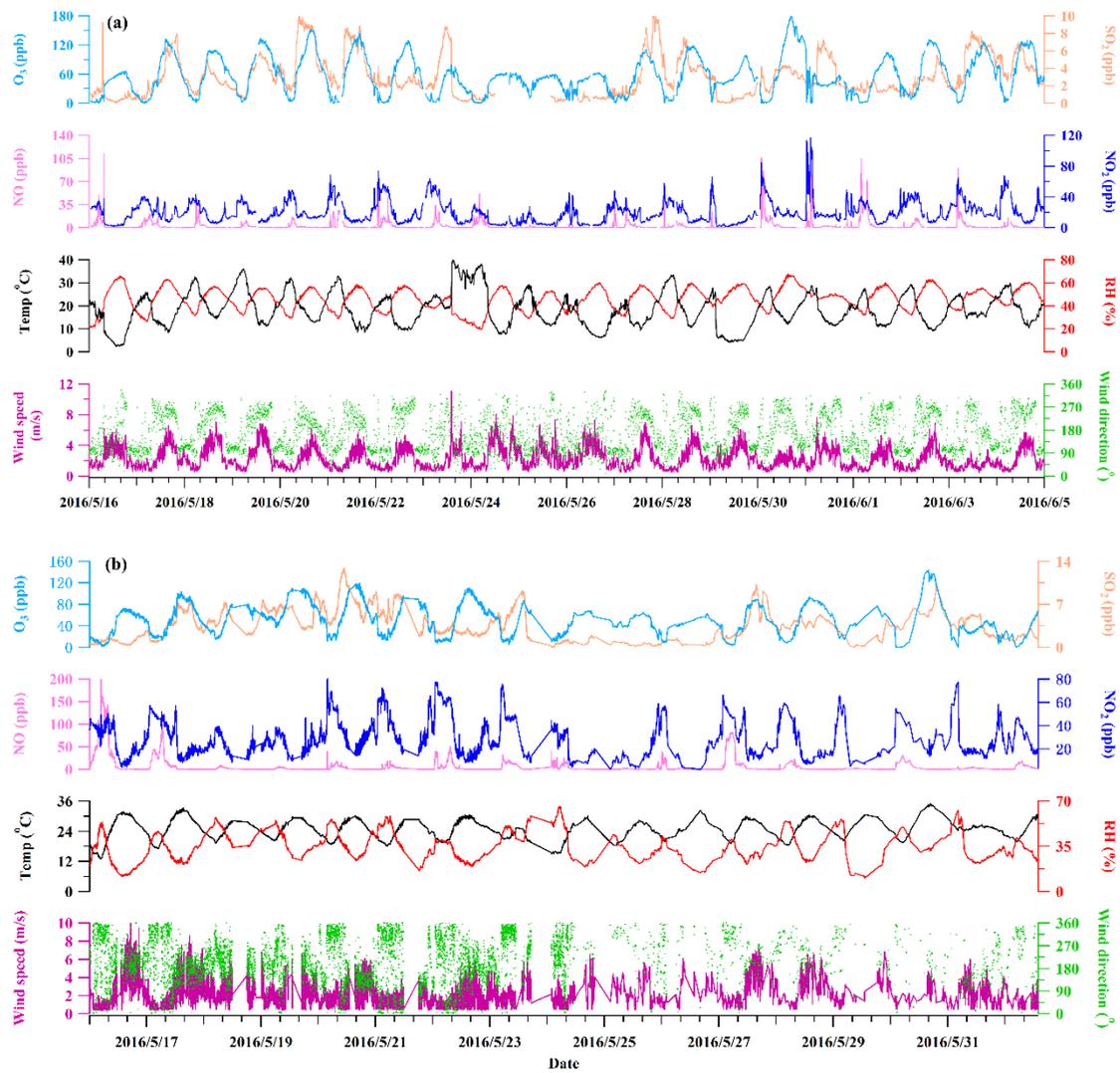


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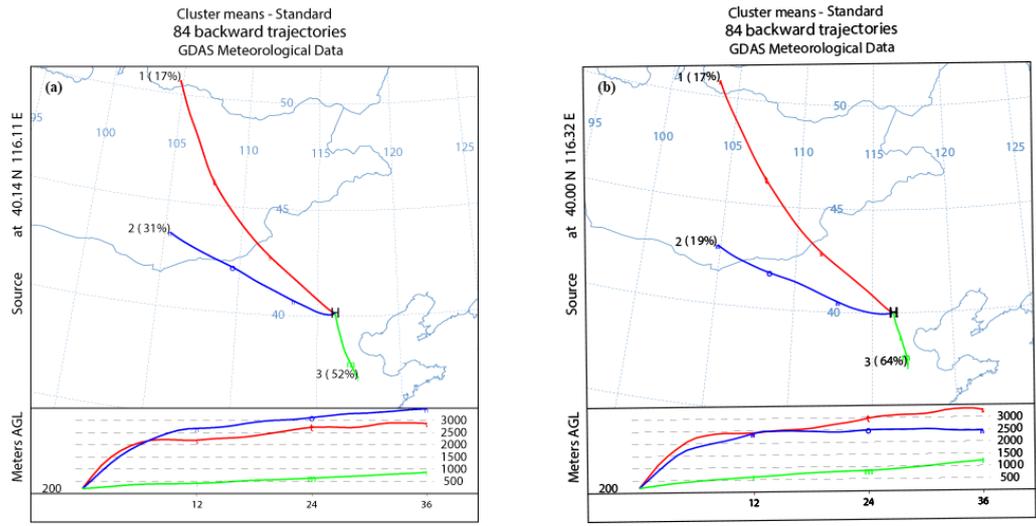


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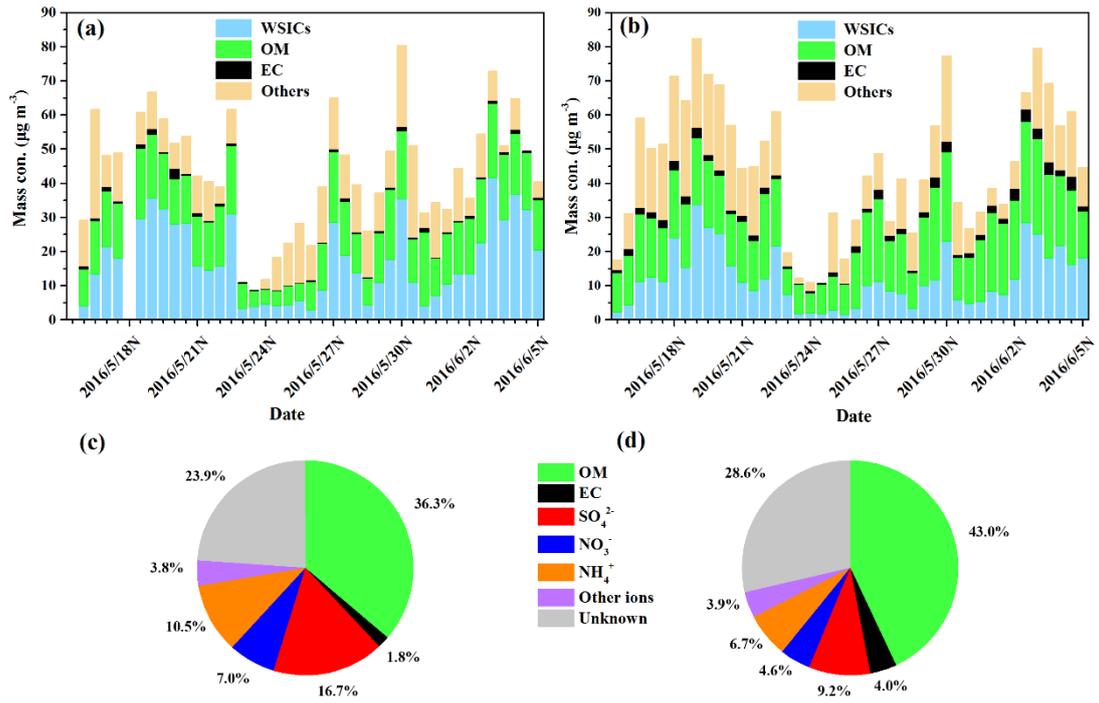


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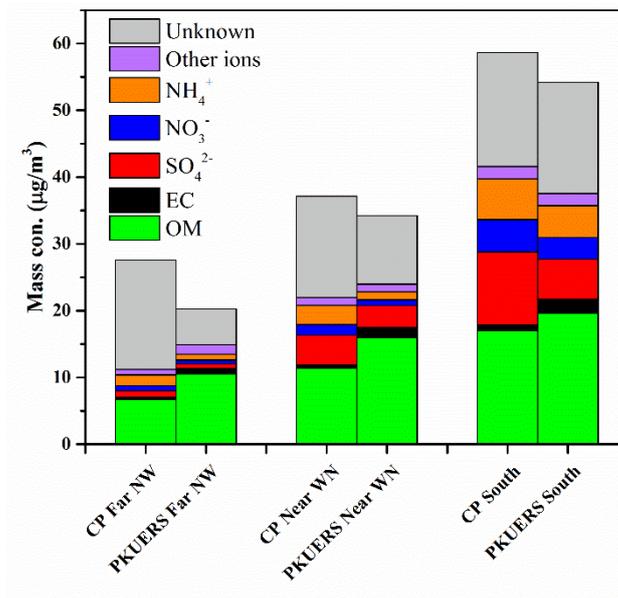


Fig. S5 Three categories of PM_{2.5} chemical compositions according to back trajectory clustering analysis

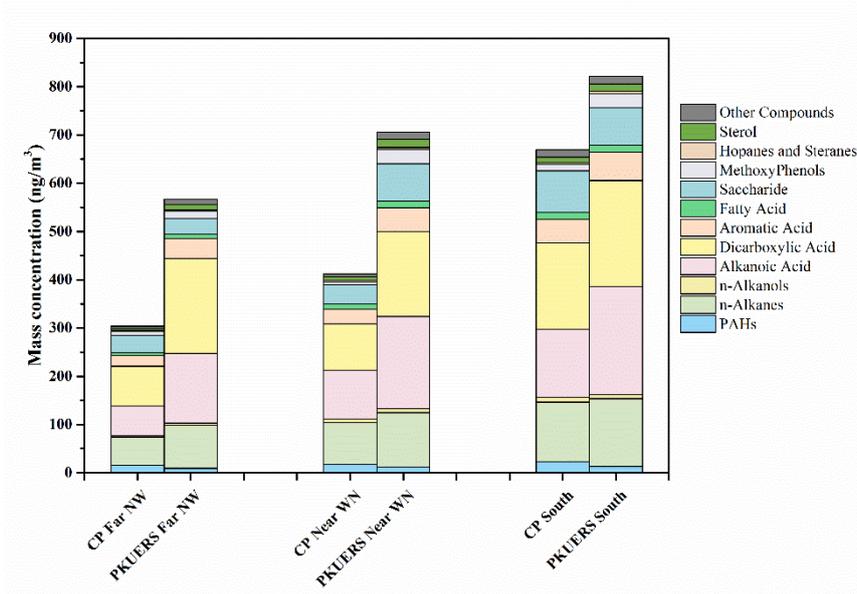


Fig. S6 Primary organic matter from different directions of CP and PKUERS

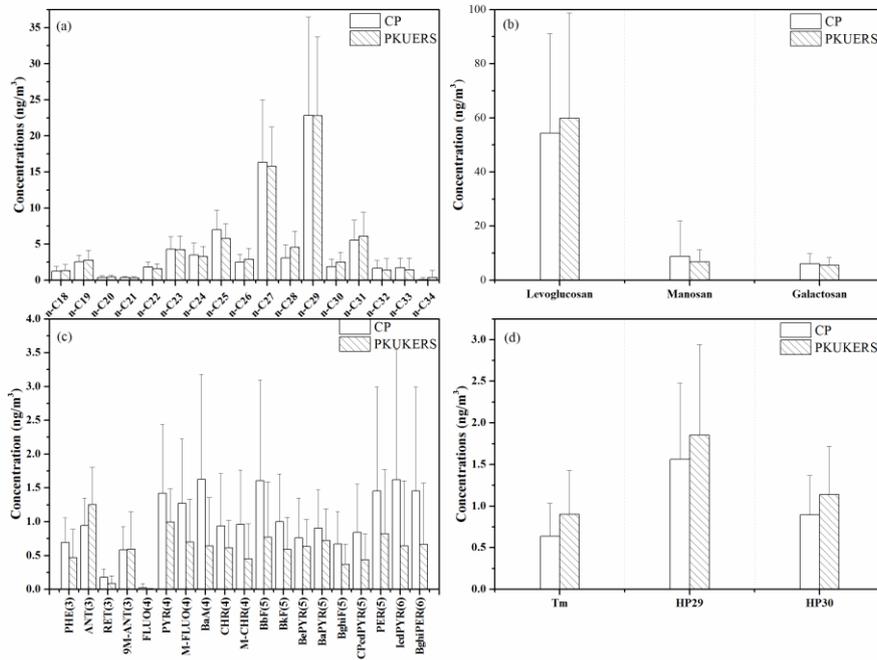


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TableTable S1 Mixing ratios of gaseous pollutants (SO₂, O₃, NO, NO₂)

Type	Sites	NO	NO ₂	O ₃	SO ₂	CO
Urban	PKUERS (2010) (Zheng et al., 2016a)	7.1	26.0	32.0	4.9	820.0
	PKUERS (2016)	10.5	32.3	48.6	3.9	623.6
	CP(2016)	3.1	19.7	57.7	3.0	492.4

Table S2 Concentrations of tracer compounds for primary sources

	CP			PKUERS		
	Daytime	Nighttime	Diurnal	Daytime	Nighttime	Diurnal
Tracers (ng m⁻³)						
n-alkanes (C28-C33)						
C-28	2.74±1.18	3.34±2.29	3.10±1.84	3.99±1.93	5.06±2.44	4.54±2.24
C-29	23.12±14.38	22.54±13.38	22.82±13.68	20.49±8.48	24.96±12.69	22.79±10.94
C-30	1.54±0.62	2.19±1.25	1.88±1.04	2.06±0.72	3.25±1.30	2.72±1.23
C-31	5.22±1.98	5.85±3.40	5.54±2.78	5.23±2.31	7.24±3.72	6.23±3.22
C-32	1.67±1.05	1.72±1.11	1.70±1.07	1.91±0.92	2.69±1.71	2.34±1.43
C-33	1.83±0.94	2.49±1.25	2.19±1.15	2.54±1.12	2.74±1.51	2.65±1.32
Levogluconan						
Levogluconan	28.86±16.95	75.92±40.91	53.03±39.26	39.09±25.04	79.56±39.97	59.87±38.93
Hopanes						
17α(H)-22,29,30-trishopane	0.61±0.28	0.65±0.44	0.63±0.36	0.88±0.55	0.93±0.52	0.90±0.53
17β(H)-21α(H)-norhopane	1.42±0.48	1.66±0.97	1.55±0.70	1.93±1.24	1.79±0.96	1.86±1.09
17α(H)-21β(H)-hopane	0.80±0.32	0.95±0.59	0.88±0.47	1.12±0.64	1.16±0.55	1.14±0.58
PAHs						
benzo(b)fluoranthene	0.64±0.45	2.45±1.68	1.57±1.53	0.53±0.47	1.07±1.00	0.81±0.82
benzo(k)fluoranthene	0.50±0.25	1.42±0.73	0.98±0.72	0.50±0.38	0.69±0.54	0.60±0.47
benzo(e)pyrene	0.38±0.22	1.09±0.67	0.74±0.61	0.59±0.39	0.70±0.40	0.65±0.39
benzo(ghi)perylene	0.71±0.40	2.48±1.63	1.81±1.56	1.15±0.43	1.84±0.80	1.54±0.73
indeno(1,2,3-cd)pyrene	0.76±0.52	2.78±2.24	2.01±2.03	0.79±0.26	2.02±0.97	1.48±0.96

Table S3 Comparison of SOA tracers for the biogenic and anthropogenic sources in different regions

Species	Tracer (ng m ⁻³)	IITB	CH	YL	CL	RTP	Yufa 2008	PKUERS 2008	PKUERS 2016	CP 2016
Isoprene	I-1	0.4±0.4	0.3±0.2	1.0±1.3	-	26.5±20.9	29.9±19.9	21.0±19.5	1.2±0.7	1.4±1.7
	I-2	-	0.2±0.1	8.3±12.9	11.2±5.8	63.4±19.2	63.6±46.8	42.2±28.2	7.7±4.7	9.9±4.6
	I-3	-	0.5±0.3	20.3±20.3	24.1±13.4	85.7±27.6	121.5±101.4	77.2±60.2	16.8±6.3	10.5±5.7
	∑isoprene	1.9±2.0	1.0±0.6	29.6±34.5	35.3±19.2	175.6±49.7	215.0±160.3	140.4±100.9	25.7±11.7	18.8±12.0
α-pinene	A-1	-	-	8.4±10.7	17.2±6.3	1.8±0.7	9.6±7.2	8.7±7.1	7.0±2.7	9.2±5.5
	A-2	-	-	5.9±4.9	17.1±5.4	16.3±7.3	5.3±2.8	6.8±7.1	4.8±2.8	4.2±2.4
	A-3	-	-	-	-	25.9±13.9	4.3±6.9	4.6±4.0	15.6±8.3	11.8±6.3
	A-4	-	-	4.6±3.6	-	46.1±18.9	7.7±5.5	8.3±5.8	4.8±1.2	4.4±1.3
	A-5	2.0±1.2	-	10.2±7.0	-	53.5±17.9	57.5±27.4	51.7±31.3	16.8±6.3	10.5±5.7
	A-6	-	-	9.7±15.1	-	12.0±1.0	8.3±6.7	9.9±9.1	13.0±4.8	13.6±8.0
	A-7	0.3±0.2	-	8.0±7.5	2.2±1.0	4.7±1.3	6.1±5.5	8.5±10.1	12.4±6.0	13.6±8.0
	PA	0.6±0.3	0.2±0.5	-	4.4±3.5	9.2±3.6	3.4±5.7	3.2±6.4	10.0±4.5	9.5±12.7
	PNA	5.5±3.9	0.1±0.1	-	6.7±1.9	-	21.9±9.2	11.9±6.3	11.9±3.3	15.4±8.1
	∑α-pinene	8.4±5.6	0.3±0.6	46.8±56.3	47.6±18.1	169.5±64.6	124.1±47.8	113.5±63.3	96.4±39.9	92.2±58.0
β-caryophyllene	C-1	0.7±0.3	0.2±0.5	12.5±16.9	1.4±0.4	25.3±4.2	3.6±2.3	5.1±4.5	6.0±2.8	6.1±3.5
Toluene	T-3	0.1±0.1	-	1.7±1.8	8.3±2.8	4.1±1.9	11.7±6.9	13.3±7.7	11.0±3.7	9.7±7.3

I-1 represented 2-methylglyceric acid, I-2 represented 2-Methylthreitol, I-3 represented 2-methylerythritol, A-1 represented 3-Isopropylpentanedioic acid, A-2 represented 3-Acetylpentanedioic acid, A-3 represented 2-Hydroxy-4-isopropyladipic acid, A-4 represented 3-Acetyl hexanedioic acid, A-5 represented 3-Hydroxyglutaric acid, A-6 represented 2-Hydroxy-4,4-dimethylglutaric acid, A-7 represented 3-(2-Hydroxy-ethyl)-2,2-dimethyl-cyclobutane-carboxylic acid, PA represented pinic acid, PNA represented pinonic acid, T-3 represented 2,3-dihydroxy-4-oxopentanoic acid, C-1 represented β-caryophyllinic acid.