

## Application Data Sheet



GC-MS

Analysis of Brominated Flame Retardants and Phthalate Esters Under the Same Conditions Using a Pyrolysis GC-MS System (4) - TBBPA and BPBPE -

LAAN-J-MS-E067

In recent years, an analysis method has been required to determine not only polybrominated biphenyls (PBBs) and polybrominated diphenyl ethers (PBDEs), which are regulated under the RoHS Directive, but also phthalate esters and other brominated flame retardants not governed by the directive (such as tetrabromobisphenol A (TBBPA), hexabromocyclododecane (HBCDD), and bis(pentabromophenyl)ethane (BPBPE). TBBPA and BPBPE are not regulated by the RoHS Directive; however, they are frequently detected as brominated flame retardant additives by the EDX screening method. This Application Data Sheet shows the results from analyzing TBBPA and BPBPE in polymers under the same analytical conditions as those in Application Data Sheet 47 using the EGA/PY-3030D Multi-Shot Pyrolyzer and GCMS-QP2020 Ultra systems

## Experimental

TBBPA and BPBPE were dissolved with toluene at a concentration of 100  $\mu$ g/mL, respectively. Polystyrene was dissolved in a mixture of 9:1 (v/v) dichloromethane and xylene at a concentration of 25 mg/mL. 20 $\mu$ L of polystyrene solution (0.5 mg) and 5  $\mu$ L of each TBBPA and BPBPE solution (0.5  $\mu$ g) were added to Eco-Cup LF. of the pyrolyzer. The solvent was evaporated to dryness at room temperature. The concentration of TBBPA and BPBPE was 1000 ppm in polystyrene. FASST (Fast Automated Scan/SIM Type), which is capable of simultaneous Scan and SIM measurements, was used as the measurement mode. Table 1 shows the analytical conditions and Fig. 1 shows the SIM measurement program.

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Pyrolysis Instrument GC-MS Column	:EGA/PY-3030D Multi-Shot Pyrolyzer :GCMS-QP2010 Ultra :Ultra ALLOY-PBDE [15 m length, 0.25 mm I.D. , df = 0.05 μm]		
[Pyrolyzer] Pyrolysis Furnace Temp. Interface Temp. [GC] Injection Temp. Column Oven Temp. Injection Mode Carrier Gas Flow Control Mode Purge Flow Rate Split Ratio	:200 °C $\rightarrow$ (20 °C/min) $\rightarrow$ 300 °C $\rightarrow$ (5 °C / :Manual (300 °C) :320 °C :80 °C $\rightarrow$ (20 °C/min) $\rightarrow$ 300 °C (5 min) :Split :Helium :Constant linear velocity (52.1 cm/sec) :3.0 mL/min :50	min) $\rightarrow$ 340 °C (1 min) [MS] Interface Temp. Ion Source Temp. Solvent Cut Time Tuning Mode Measurement Mode Scan Mass Range Scan Event Time Scan Speed SIM Monitoring m/z: SIM Event Time SIM Micro-Scan Width	:320 °C :230 °C :0.5 min :Normal :Scan/SIM : <i>m/z</i> 50 - 1000 :0.15 sec :10,000 <i>u</i> /sec See Fig. 2. :0.3 sec :0.5 <i>u</i>

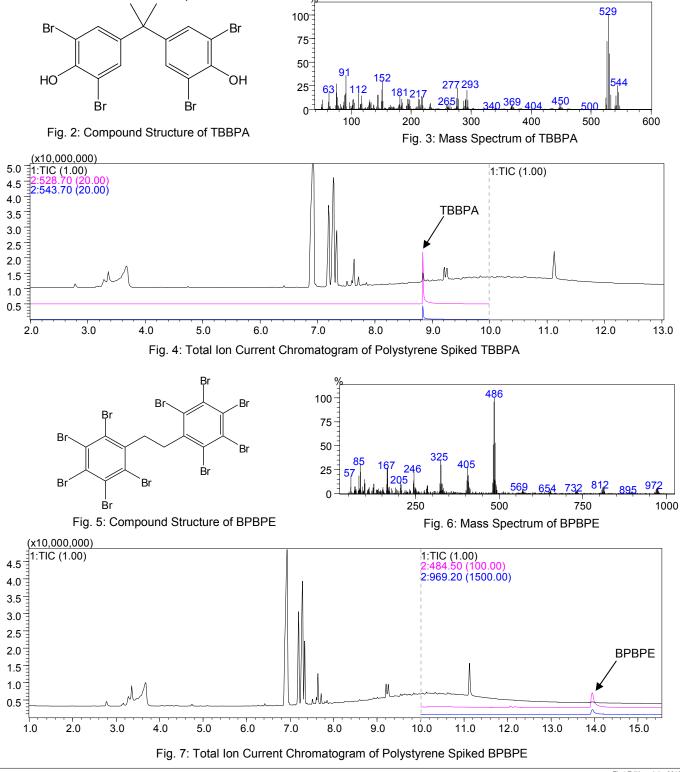
Group 1 Group 2 16 min 1 min 10 min (No. of *m/z* channels: 21) (No. of m/z channels: 11) Tetra-BDE (m/z 325.9, 483.7) Hexa-BDE (m/z 483.7, 641.5) Penta-BDE (m/z 403.8, 563.6) Hepta-BDE (m/z 563.6, 721.4) Hexa-BDE (m/z 483.7, 641.5) Octa-BDE (m/z 641.5, 801.3) Hepta-BDE (m/z 563.6, 721.4) Nona-BDE (m/z 719.4, 721.4) Tetrabromobisphenol A [TBBPA] (m/z 528.7, 543.7) Deca-BDE (m/z 799.3, 801.3) Hexabromocyclododecane [HBCDD] (m/z 319.1, 560.6) Deca-BB (m/z 941.3, 943.3) Diisobutyl phthalate [DIBP] (m/z 149.0, 205.1, 223.1) Bis(pentabromophenyl)ethane [BPBPE] Di-n-butyl phthalate [DIBP] (m/z 149.0, 205.1, 223.1) (m/z 484.5, 969.2) Benzylbutyl phthalate [BBP] (m/z 91.0, 149.0, 206.1) Bis(2-ehtylhexyl) phthalate [DEHP] (m/z 149.0, 167.0, 279.1) Di-n-octyll phthalate [DOP] (m/z 149.0, 261.1, 279.1) Di-isononyl phthalate [DINP] (m/z 149.0, 167.0, 293.1) Di-isodecyl phthalate [DIDP] (m/z 149.0, 167.0, 307.1)

## Table 1: Analytical Conditions

## Results

The chemical structure and mass spectrum of TBBPA are shown in Fig. 2 and 3. Fig. 4 shows the total ion current chromatogram of TBBPA in polystyrene (1000 ppm). The chemical structure and mass spectrum of BPBPE are shown in Fig. 5 and 6. Fig. 7 shows the total ion current chromatogram of BPBPE in polystyrene (1000 ppm).

These analytical conditions can be applied to analyses of PBBs, PBDEs ,and phthalate esters. Moreover, using the scan/ SIM mode (FASST) enable accurate determination of target compounds from the SIM data and identification of unknown compounds from the full-scan data.



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