

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU040521\  
 Data File : VU042923.D  
 Acq On : 05 Apr 2021 12:13  
 Operator : SY/MD  
 Sample : VU0405MBL01  
 Misc : 5.0g/10mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 VU0405MBL01

Quant Time: Apr 06 01:16:22 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\82U032421W.M  
 Quant Title : SW846 8260  
 QLast Update : Thu Mar 25 05:42:00 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units   | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| -----                       |        |      |          |       |         |          |
| Internal Standards          |        |      |          |       |         |          |
| 1) Pentafluorobenzene       | 5.382  | 168  | 128249   | 50.00 | ug/l    | 0.00     |
| 34) 1,4-Difluorobenzene     | 6.259  | 114  | 204633   | 50.00 | ug/l    | 0.00     |
| 63) Chlorobenzene-d5        | 9.426  | 117  | 195739   | 50.00 | ug/l    | 0.00     |
| 72) 1,4-Dichlorobenzene-d4  | 11.822 | 152  | 93913    | 50.00 | ug/l    | 0.00     |
| System Monitoring Compounds |        |      |          |       |         |          |
| 33) 1,2-Dichloroethane-d4   | 5.713  | 65   | 102533   | 50.15 | ug/l    | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | =     | 100.30% |          |
| 35) Dibromofluoromethane    | 5.298  | 113  | 72513    | 47.48 | ug/l    | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | =     | 94.96%  |          |
| 50) Toluene-d8              | 7.906  | 98   | 257558   | 47.94 | ug/l    | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | =     | 95.88%  |          |
| 62) 4-Bromofluorobenzene    | 10.642 | 95   | 92386    | 44.25 | ug/l    | 0.00     |
| Spiked Amount               | 50.000 |      | Recovery | =     | 88.50%  |          |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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