

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091621\
 Data File : VN068507.D
 Acq On : 16 Sep 2021 20:26
 Operator : JC/MD
 Sample : PB139158TB
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB139158TB

Quant Time: Sep 17 02:03:32 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090721W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 10 04:04:05 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.091	168	532287	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	779874	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	784394	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	318015	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	266701	54.622	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery	= 109.240%		
35) Dibromofluoromethane	8.024	113	242078	51.466	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery	= 102.940%		
50) Toluene-d8	10.446	98	960763	56.232	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery	= 112.460%		
62) 4-Bromofluorobenzene	12.733	95	305145	55.813	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery	= 111.620%		
Target Compounds						Qvalue
16) Acetone	4.301	43	14141	8.752	ug/l	95
20) Methylene Chloride	5.098	84	25661	4.373	ug/l	92
25) 2-Butanone	7.351	43	14188	5.980	ug/l	94
43) Isopropyl Acetate	8.560	43	40211	4.608	ug/l #	88

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091621\
Data File : VN068507.D
Acq On : 16 Sep 2021 20:26
Operator : JC/MD
Sample : PB139158TB
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB139158TB

Quant Time: Sep 17 02:03:32 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090721W.M
Quant Title : SW846 8260
QLast Update : Fri Sep 10 04:04:05 2021
Response via : Initial Calibration

