

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081321\
 Data File : VX023699.D
 Acq On : 13 Aug 2021 18:34
 Operator : JC/MD
 Sample : PB138358ZHE#16
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB138358ZHE#16

Quant Time: Aug 14 05:29:05 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 23 06:05:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.568	168	147559	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	256698	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	244557	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	99279	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	105571	57.355	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 114.700%			
35) Dibromofluoromethane	5.397	113	84085	52.390	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 104.780%			
50) Toluene-d8	8.653	98	321872	54.584	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 109.160%			
62) 4-Bromofluorobenzene	11.085	95	109576	54.647	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 109.300%			
Target Compounds					Qvalue	
16) Acetone	2.386	43	27068	29.978	ug/l	98
20) Methylene Chloride	2.794	84	11417	6.500	ug/l #	89
43) Isopropyl Acetate	6.348	43	22813	5.028	ug/l #	90

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081321\
Data File : VX023699.D
Acq On : 13 Aug 2021 18:34
Operator : JC/MD
Sample : PB138358ZHE#16
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB138358ZHE#16

Quant Time: Aug 14 05:29:05 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
Quant Title : SW846 8260
QLast Update : Fri Jul 23 06:05:38 2021
Response via : Initial Calibration

