

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042021\
 Data File : VX021769.D
 Acq On : 20 Apr 2021 21:25
 Operator : JC/MD
 Sample : PB135856ZHE#03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB135856ZHE#03

Quant Time: Apr 21 11:25:23 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042021W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 20 13:46:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	157120	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.830	114	262267	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.092	117	228746	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.061	152	100030	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.031	65	89269	50.08	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 100.16%			
35) Dibromofluoromethane	5.464	113	84245	50.23	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.46%			
50) Toluene-d8	8.695	98	311137	48.25	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 96.50%			
62) 4-Bromofluorobenzene	11.116	95	106502	44.22	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 88.44%			
Target Compounds						Qvalue
3) Chloromethane	1.319	50	3485	2.05	ug/l	99
16) Acetone	2.422	43	24292	39.64	ug/l	100
20) Methylene Chloride	2.837	84	10619	5.78	ug/l	95
25) 2-Butanone	4.635	43	10578	10.19	ug/l	99
43) Isopropyl Acetate	6.409	43	18027	4.73	ug/l	99

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

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