

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042021\
 Data File : VX021770.D
 Acq On : 20 Apr 2021 21:49
 Operator : JC/MD
 Sample : PB135856ZHE#04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB135856ZHE#04

Quant Time: Apr 21 11:25:42 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	159937	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.830	114	265459	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.098	117	228664	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.061	152	96740	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.031	65	89538	49.35	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.70%
35) Dibromofluoromethane	5.464	113	84303	49.66	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.32%
50) Toluene-d8	8.696	98	312421	47.86	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.72%
62) 4-Bromofluorobenzene	11.122	95	104081	42.70	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	85.40%
Target Compounds						
					Qvalue	
3) Chloromethane	1.319	50	3270	1.91	ug/l	91
16) Acetone	2.422	43	25326	40.60	ug/l	98
20) Methylene Chloride	2.837	84	10528	5.63	ug/l	96
25) 2-Butanone	4.641	43	10444	9.89	ug/l	92
43) Isopropyl Acetate	6.415	43	18082	4.69	ug/l	99

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

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