

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042021\
 Data File : VX021767.D
 Acq On : 20 Apr 2021 20:39
 Operator : JC/MD
 Sample : PB135856ZHE#01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB135856ZHE#01

Quant Time: Apr 21 11:24:37 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	167508	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.830	114	275720	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.092	117	241824	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.061	152	103337	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.031	65	94563	49.76	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.52%
35) Dibromofluoromethane	5.465	113	87360	49.55	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.10%
50) Toluene-d8	8.696	98	326519	48.16	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.32%
62) 4-Bromofluorobenzene	11.116	95	108581	42.89	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	85.78%
Target Compounds						
					Qvalue	
3) Chloromethane	1.319	50	3911	2.14	ug/l	93
16) Acetone	2.422	43	25152	38.50	ug/l	98
20) Methylene Chloride	2.837	84	10773	5.50	ug/l	90
25) 2-Butanone	4.648	43	10434	9.43	ug/l #	88
43) Isopropyl Acetate	6.409	43	18616	4.65	ug/l	97

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

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