

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060222\
 Data File : VU048825.D
 Acq On : 03 Jun 2022 00:30
 Operator : SY/MD
 Sample : N3119-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 40016

Quant Time: Jun 03 04:03:50 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 27 06:42:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.372	168	328510	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.247	114	555288	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	502288	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	249894	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	228841	46.384	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	92.760%
35) Dibromofluoromethane	5.289	113	176951	44.667	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.340%
50) Toluene-d8	7.896	98	670646	47.226	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	94.460%
62) 4-Bromofluorobenzene	10.632	95	247001	45.991	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	91.980%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	3.179	59	12803	10.236	ug/l #	38
16) Acetone	2.623	43	137200	54.568	ug/l	100
25) 2-Butanone	4.703	43	44326	12.272	ug/l	96
29) Tetrahydrofuran	5.057	42	13753	6.045	ug/l	97
51) 4-Methyl-2-Pentanone	7.790	43	69097	10.031	ug/l	97

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

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