

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060722\
 Data File : VU048917.D
 Acq On : 07 Jun 2022 18:11
 Operator : SY/MD
 Sample : N3165-19
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 43BR-20220601

Quant Time: Jun 08 04:55:09 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	237407	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	385537	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	356263	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	156429	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	180341	50.581	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 101.160%	
35) Dibromofluoromethane	5.288	113	141238	51.350	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.700%	
50) Toluene-d8	7.899	98	471496	47.821	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 95.640%	
62) 4-Bromofluorobenzene	10.632	95	163905	43.956	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 87.920%	
Target Compounds						
						Qvalue
16) Acetone	2.629	43	7749	4.265	ug/l	96
37) Ethyl Acetate	4.803	43	121814	24.004	ug/l	99
40) Benzene	5.774	78	3921	0.331	ug/l	93

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

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