

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060722\
 Data File : VU048922.D
 Acq On : 07 Jun 2022 20:09
 Operator : SY/MD
 Sample : N3166-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 55BR-20220601

Quant Time: Jun 08 04:56:41 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	248989	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	395513	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	365731	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	164236	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	182510	48.808	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.620%
35) Dibromofluoromethane	5.288	113	141734	50.230	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.460%
50) Toluene-d8	7.899	98	483820	47.833	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.660%
62) 4-Bromofluorobenzene	10.632	95	172421	45.074	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.140%
Target Compounds						
					Qvalue	
16) Acetone	2.626	43	7160	3.757	ug/l #	87
37) Ethyl Acetate	4.803	43	191442	36.772	ug/l	98
52) Toluene	7.973	92	3320	0.457	ug/l	87

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

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