

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060722\
 Data File : VU048936.D
 Acq On : 08 Jun 2022 02:14
 Operator : SY/MD
 Sample : N3166-09
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DUP06-20220531

Quant Time: Jun 08 09:15:21 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	247688	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	404040	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	376578	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	170240	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	188781	50.750	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.500%
35) Dibromofluoromethane	5.288	113	142085	49.292	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.580%
50) Toluene-d8	7.896	98	502613	48.643	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.280%
62) 4-Bromofluorobenzene	10.632	95	183414	46.936	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.880%
Target Compounds						Qvalue
16) Acetone	2.626	43	5953	3.140	ug/l	97
27) cis-1,2-Dichloroethene	4.665	96	1794	0.537	ug/l	51
37) Ethyl Acetate	4.803	43	109504	20.590	ug/l	98

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

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