

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060722\
 Data File : VU048932.D
 Acq On : 08 Jun 2022 00:39
 Operator : SY/MD
 Sample : N3164-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 28BR-20220531

Quant Time: Jun 08 09:14:05 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	245623	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	405537	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	364458	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	149695	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	177724	48.180	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.360%
35) Dibromofluoromethane	5.288	113	138824	47.983	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.960%
50) Toluene-d8	7.899	98	488218	47.075	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	94.160%
62) 4-Bromofluorobenzene	10.632	95	166193	42.372	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	84.740%
Target Compounds						
					Qvalue	
16) Acetone	2.623	43	5450	2.899	ug/l #	83
30) Chloroform	5.089	83	2605	0.464	ug/l	86
37) Ethyl Acetate	4.800	43	150524	28.198	ug/l	100

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

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