

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063022\
 Data File : VU049567.D
 Acq On : 30 Jun 2022 16:44
 Operator : SY/MD
 Sample : N3543-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 KENS-062822-SB18-GW

Quant Time: Jul 01 05:26:17 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 28 10:35:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.372	168	235206	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.250	114	409471	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	382680	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	180115	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	166519	45.970	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	91.940%
35) Dibromofluoromethane	5.288	113	130159	45.435	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.860%
50) Toluene-d8	7.896	98	499992	48.933	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.860%
62) 4-Bromofluorobenzene	10.632	95	185108	47.292	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.580%
Target Compounds						
					Qvalue	
16) Acetone	2.626	43	4332	2.371	ug/l	# 86
17) Carbon Disulfide	2.787	76	2490	0.313	ug/l	# 86
30) Chloroform	5.089	83	7176	1.281	ug/l	97
40) Benzene	5.777	78	4777	0.381	ug/l	95
52) Toluene	7.970	92	5464	0.733	ug/l	94
95) Naphthalene	14.089	128	24692	3.426	ug/l	97

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

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