

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
 Data File : VX029354.D
 Acq On : 09 Jun 2022 18:45
 Operator : JC/MD
 Sample : N3266-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OH-WATER-0608

Quant Time: Jun 10 08:10:18 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 08 03:00:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	200225	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	327402	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	308278	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	165714	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	177664	58.593	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 117.180%	
35) Dibromofluoromethane	5.391	113	159040	57.139	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 114.280%	
50) Toluene-d8	8.653	98	498010	50.590	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 101.180%	
62) 4-Bromofluorobenzene	11.085	95	231734	60.154	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 120.300%	
Target Compounds						
					Qvalue	
5) Bromomethane	1.617	94	1031	0.547	ug/l	97
16) Acetone	2.386	43	14771	9.119	ug/l	97
17) Carbon Disulfide	2.514	76	2268	0.306	ug/l	96
25) 2-Butanone	4.580	43	4315	1.684	ug/l	96
39) Methylcyclohexane	7.385	83	2235	0.395	ug/l	90
52) Toluene	8.726	92	3991	0.454	ug/l	84

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

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