

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112224\
 Data File : VX043955.D
 Acq On : 22 Nov 2024 13:25
 Operator : JC/MD
 Sample : P4947-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 A3988

Quant Time: Nov 22 22:21:57 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 22 03:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	113610	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	217300	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	188960	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	77731	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	98276	50.434	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	100.860%
35) Dibromofluoromethane	5.379	113	72758	46.858	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.720%
50) Toluene-d8	8.647	98	262229	48.993	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	97.980%
62) 4-Bromofluorobenzene	11.079	95	87848	48.227	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	96.460%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	18996	21.225	ug/l	99
17) Carbon Disulfide	2.502	76	1018	0.377	ug/l	99
20) Methylene Chloride	2.782	84	437	0.288	ug/l #	84
68) m/p-Xylenes	10.305	106	1172	0.447	ug/l	79
69) o-Xylene	10.640	106	951	0.371	ug/l	97

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

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