

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112524\
 Data File : VX043993.D
 Acq On : 25 Nov 2024 17:44
 Operator : JC/MD
 Sample : P4873-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-MW-28-20241114

Quant Time: Nov 26 00:51:39 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	5.550	168	108714	50.000	ug/l	0.01
34) 1,4-Difluorobenzene	6.757	114	206635	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	183586	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	69483	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	94974	50.934	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	101.860%	
35) Dibromofluoromethane	5.379	113	69574	47.121	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	94.240%	
50) Toluene-d8	8.647	98	259104	50.908	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	101.820%	
62) 4-Bromofluorobenzene	11.079	95	79414	45.847	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	91.700%	
Target Compounds						
16) Acetone	2.386	43	2595	3.030	ug/l	93
20) Methylene Chloride	2.782	84	509	0.350	ug/l #	77
21) trans-1,2-Dichloroethene	3.081	96	1036	0.801	ug/l #	72
30) Chloroform	5.093	83	3767	1.387	ug/l #	79
32) 1,1,1-Trichloroethane	5.373	97	8101	3.459	ug/l #	87
44) Trichloroethene	7.117	130	405015	249.635	ug/l	94
47) Bromodichloromethane	7.824	83	3405	1.710	ug/l #	94
60) Dibromochloromethane	9.525	129	2389	1.682	ug/l	96
64) Tetrachloroethene	9.269	164	29576	24.432	ug/l	96

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

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