

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112524\
 Data File : VX043994.D
 Acq On : 25 Nov 2024 18:08
 Operator : JC/MD
 Sample : P4873-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-MW-29-20241114

Quant Time: Nov 26 00:52:06 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	112896	50.000	ug/l	0.01
34) 1,4-Difluorobenzene	6.757	114	215968	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	189479	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	83328	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	97434	50.318	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 100.640%			
35) Dibromofluoromethane	5.385	113	73436	47.587	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 95.180%			
50) Toluene-d8	8.647	98	262792	49.401	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 98.800%			
62) 4-Bromofluorobenzene	11.079	95	88246	48.744	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 97.480%			
Target Compounds						
16) Acetone	2.386	43	12205	13.723	ug/l	97
20) Methylene Chloride	2.782	84	624	0.413	ug/l #	80
21) trans-1,2-Dichloroethene	3.081	96	518	0.386	ug/l	90
27) cis-1,2-Dichloroethene	4.489	96	5547	3.341	ug/l	91
30) Chloroform	5.099	83	2172	0.770	ug/l	92
32) 1,1,1-Trichloroethane	5.385	97	7204	2.962	ug/l	99
44) Trichloroethene	7.123	130	233374	137.626	ug/l	92
47) Bromodichloromethane	7.824	83	1603	0.770	ug/l #	85
60) Dibromochloromethane	9.525	129	1084	0.730	ug/l	89
64) Tetrachloroethene	9.275	164	13688	10.956	ug/l	96

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

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