

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120924\
 Data File : VX044190.D
 Acq On : 09 Dec 2024 16:04
 Operator : JC/MD
 Sample : P5145-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 286085

Quant Time: Dec 10 01:44:24 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	105331	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	198577	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	173539	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	76204	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	91293	50.533	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 101.060%			
35) Dibromofluoromethane	5.379	113	64266	45.292	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 90.580%			
50) Toluene-d8	8.647	98	242903	49.661	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 99.320%			
62) 4-Bromofluorobenzene	11.079	95	84067	50.502	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 101.000%			
Target Compounds						
11) Tert butyl alcohol	2.983	59	2406	8.297	ug/l #	95
16) Acetone	2.392	43	16875	20.337	ug/l	97
18) Methyl Acetate	2.709	43	125584	64.992	ug/l	99
25) 2-Butanone	4.587	43	10061	9.397	ug/l	97
29) Tetrahydrofuran	5.044	42	2915	4.441	ug/l #	89
51) 4-Methyl-2-Pentanone	8.586	43	1110	0.521	ug/l	99

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

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