

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022124\
 Data File : VX040389.D
 Acq On : 21 Feb 2024 18:08
 Operator : JC/MD
 Sample : P1532-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE104D2-20240216

Quant Time: Feb 22 01:12:23 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
 Quant Title : SW846 8260
 QLast Update : Mon Feb 12 00:21:14 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	72031	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	130211	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	124358	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	60248	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	57453	53.614	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 107.220%			
35) Dibromofluoromethane	5.391	113	43434	47.765	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 95.540%			
50) Toluene-d8	8.647	98	158214	46.513	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 93.020%			
62) 4-Bromofluorobenzene	11.079	95	59829	43.127	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 86.260%			
Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.319	96	1554	2.113	ug/l #	91
16) Acetone	2.374	43	1564	3.554	ug/l	97
21) trans-1,2-Dichloroethene	3.099	96	341	0.407	ug/l #	63
24) 1,1-Dichloroethane	3.611	63	3206	2.126	ug/l #	93
27) cis-1,2-Dichloroethene	4.489	96	53327	55.812	ug/l	90
30) Chloroform	5.105	83	3705	2.273	ug/l	90
42) 1,2-Dichloroethane	6.086	62	3196	2.295	ug/l	98
44) Trichloroethene	7.123	130	433674	441.360	ug/l	96
52) Toluene	8.720	92	9759	4.183	ug/l	93
64) Tetrachloroethene	9.275	164	507	0.582	ug/l #	77

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

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