

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010324\
 Data File : VX039594.D
 Acq On : 03 Jan 2024 12:12
 Operator : JC/MD
 Sample : P1001-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ORA-1206

Quant Time: Jan 04 03:29:33 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 14 01:28:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	94837	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	198441	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	201316	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	87104	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	82858	47.692	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	95.380%		
35) Dibromofluoromethane	5.385	113	65486	47.768	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.540%		
50) Toluene-d8	8.647	98	250537	49.225	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.460%		
62) 4-Bromofluorobenzene	11.079	95	105181	50.087	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	100.180%		
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	3.001	59	6316	14.477	ug/l #	90
16) Acetone	2.385	43	44514	22.828	ug/l	93
18) Methyl Acetate	2.703	43	6225	1.657	ug/l	97
20) Methylene Chloride	2.782	84	565	0.390	ug/l #	75
25) 2-Butanone	4.562	43	34454	22.982	ug/l	94
51) 4-Methyl-2-Pentanone	8.586	43	1989	0.718	ug/l	97
52) Toluene	8.720	92	774	0.210	ug/l	76
59) 2-Hexanone	9.451	43	1321	0.553	ug/l	89
68) m/p-Xylenes	10.299	106	763	0.260	ug/l	89
84) 1,2,4-Trimethylbenzene	11.756	105	1494	0.254	ug/l	93
95) Naphthalene	13.780	128	3499	0.549	ug/l	97

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

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