

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012725\
 Data File : VX044719.D
 Acq On : 27 Jan 2025 13:24
 Operator : JC/MD
 Sample : Q1173-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-192-GW-100-102

Quant Time: Jan 28 04:11:28 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011525W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jan 16 01:16:09 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	156971	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	304234	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	268221	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	115943	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	113380	47.853	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	95.700%		
35) Dibromofluoromethane	5.379	113	98218	50.818	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	101.640%		
50) Toluene-d8	8.647	98	372179	55.227	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	110.460%		
62) 4-Bromofluorobenzene	11.079	95	131621	52.391	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	104.780%		
Target Compounds						
16) Acetone	2.380	43	7094	7.568	ug/l	97
30) Chloroform	5.086	83	4072	1.094	ug/l	94
44) Trichloroethene	7.135	130	1739	0.868	ug/l	74
47) Bromodichloromethane	7.818	83	3219	0.983	ug/l #	90
60) Dibromochloromethane	9.525	129	2553	1.054	ug/l	99

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

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