

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110724\
 Data File : VX043745.D
 Acq On : 07 Nov 2024 18:35
 Operator : JC/MD
 Sample : P4739-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-G2

Quant Time: Nov 08 04:13:12 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
 Quant Title : SW846 8260
 QLast Update : Mon Oct 28 13:41:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	83642	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	150865	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	132249	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	58071	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	48461	36.336	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	72.680%#	
35) Dibromofluoromethane	5.385	113	43920	42.932	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	85.860%	
50) Toluene-d8	8.647	98	172147	48.611	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	97.220%	
62) 4-Bromofluorobenzene	11.079	95	57659	43.940	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	87.880%	
Target Compounds						
						Qvalue
16) Acetone	2.404	43	9971	18.827	ug/l	94
18) Methyl Acetate	2.727	43	2103	1.359	ug/l	97
20) Methylene Chloride	2.782	84	2320	2.033	ug/l #	91
40) Benzene	6.038	78	7947	1.964	ug/l #	85
43) Isopropyl Acetate	6.348	43	46232	18.983	ug/l	97

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

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