

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
 Data File : VX043129.D
 Acq On : 24 Sep 2024 17:44
 Operator : JC/MD
 Sample : P4164-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE134D1-20240919

Quant Time: Sep 25 01:33:37 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 24 03:13:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	92513	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	182580	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	178912	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	96078	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	81271	52.887	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 105.780%			
35) Dibromofluoromethane	5.391	113	62856	48.802	ug/l	0.01
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.600%			
50) Toluene-d8	8.647	98	224968	51.628	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 103.260%			
62) 4-Bromofluorobenzene	11.079	95	95463	54.400	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 108.800%			
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	1273	1.771	ug/l #	88
30) Chloroform	5.098	83	971	0.447	ug/l #	72
40) Benzene	6.043	78	1787	0.383	ug/l	96
44) Trichloroethene	7.123	130	11707	9.162	ug/l	87
52) Toluene	8.720	92	1623	0.540	ug/l	98
64) Tetrachloroethene	9.275	164	930	0.717	ug/l #	80
68) m/p-Xylenes	10.305	106	1522	0.613	ug/l	92
69) o-Xylene	10.640	106	992	0.398	ug/l	93
84) 1,2,4-Trimethylbenzene	11.756	105	1724	0.291	ug/l	92

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

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