

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
 Data File : VX037634.D
 Acq On : 09 Sep 2023 17:50
 Operator : JC/MD
 Sample : 04297-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1731-DRIP

Quant Time: Sep 11 01:26:15 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 09 06:38:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	135558	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	221743	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	198591	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	86459	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	75089	41.035	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	82.080%		
35) Dibromofluoromethane	5.391	113	67043	43.227	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	86.460%		
50) Toluene-d8	8.647	98	260249	45.750	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	91.500%#		
62) 4-Bromofluorobenzene	11.079	95	93394	43.563	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	87.120%		
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.959	59	3125	7.632	ug/l #	93
16) Acetone	2.380	43	20772	27.604	ug/l #	88
17) Carbon Disulfide	2.514	76	6153	1.688	ug/l	98
25) 2-Butanone	4.580	43	4348	3.810	ug/l #	66
51) 4-Methyl-2-Pentanone	8.580	43	2140	1.069	ug/l	100
68) m/p-Xylenes	10.305	106	4495	1.541	ug/l	97
69) o-Xylene	10.646	106	1349	0.465	ug/l	98
80) 1,3,5-Trimethylbenzene	11.451	105	3469	0.630	ug/l	98
84) 1,2,4-Trimethylbenzene	11.756	105	11617	2.109	ug/l	94
95) Naphthalene	13.780	128	3522	0.547	ug/l	97

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

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