

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121223\
 Data File : VX039172.D
 Acq On : 12 Dec 2023 17:43
 Operator : JC/MD
 Sample : 05757-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OF-2

Quant Time: Dec 13 04:19:04 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111623W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 16 22:59:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	109642	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	226662	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	231340	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	116913	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	95039	53.794	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 107.580%	
35) Dibromofluoromethane	5.385	113	74119	45.277	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 90.560%	
50) Toluene-d8	8.647	98	282848	45.160	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 90.320%#	
62) 4-Bromofluorobenzene	11.079	95	129465	51.637	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 103.280%	
Target Compounds						
					Qvalue	
16) Acetone	2.374	43	22566	23.994	ug/l	91
19) Methyl tert-butyl Ether	3.111	73	2127	0.443	ug/l	92
31) Cyclohexane	5.471	56	586	0.232	ug/l #	68
39) Methylcyclohexane	7.373	83	695	0.227	ug/l #	84
51) 4-Methyl-2-Pentanone	8.574	43	18838	6.034	ug/l	99
65) Chlorobenzene	10.080	112	6171	1.238	ug/l	95

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

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