

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031623\
 Data File : VX034571.D
 Acq On : 17 Mar 2023 07:59
 Operator : JC/MD
 Sample : 01907-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 33959

Quant Time: Mar 17 08:34:48 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Mar 13 10:29:32 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	261696	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	521941	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	418382	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	198901	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	229335	60.238	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 120.480%#	
35) Dibromofluoromethane	5.385	113	143318	41.588	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 83.180%	
50) Toluene-d8	8.647	98	633647	49.255	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 98.520%	
62) 4-Bromofluorobenzene	11.079	95	225337	43.365	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 86.740%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.953	59	13296	16.771	ug/l #	93
16) Acetone	2.374	43	135292	75.441	ug/l	99
17) Carbon Disulfide	2.514	76	18359	2.277	ug/l	98
18) Methyl Acetate	2.709	43	7822	1.335	ug/l	99
20) Methylene Chloride	2.788	84	13081	3.758	ug/l	98
25) 2-Butanone	4.562	43	21968	8.561	ug/l	100
37) Ethyl Acetate	4.715	43	61360	10.611	ug/l	97
40) Benzene	6.038	78	47095	3.042	ug/l	99
43) Isopropyl Acetate	6.336	43	284413	29.034	ug/l	99
51) 4-Methyl-2-Pentanone	8.574	43	540581	95.128	ug/l #	79
52) Toluene	8.720	92	31897	3.327	ug/l	99
68) m/p-Xylenes	10.299	106	21332	3.379	ug/l	97
69) o-Xylene	10.640	106	30371	4.794	ug/l	99
95) Naphthalene	13.774	128	249756	17.054	ug/l	99

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

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