

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112120\
 Data File : VX019521.D
 Acq On : 21 Nov 2020 02:47
 Operator : JC/MD
 Sample : L4779-14 10X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 COMP-1605-1607

Manual Integrations
 APPROVED

MMDadoda
 11/23/2020 11:05:59 AM

Quant Time: Nov 21 06:22:15 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 21 01:57:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.62	168	281787	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	461953	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	405771	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	195035	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	6.03	65	173358	47.45	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	94.90%
35) Dibromofluoromethane	5.46	113	135582	46.76	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	93.52%
50) Toluene-d8	8.70	98	545443	48.41	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.82%
62) 4-Bromofluorobenzene	11.12	95	196760	48.53	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	97.06%

Target Compounds

					Qvalue
11) Tert butyl alcohol	3.00	59	10830m	15.663	ug/l
16) Acetone	2.44	43	22817	16.149	ug/l
17) Carbon Disulfide	2.53	76	8427	1.121	ug/l #
18) Methyl Acetate	2.76	43	1909	1.179	ug/l
25) 2-Butanone	4.66	43	5386	2.538	ug/l
51) 4-Methyl-2-Pentanone	8.63	43	7694	1.931	ug/l
84) 1,2,4-Trimethylbenzene	11.79	105	2860	0.250	ug/l

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

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