

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042221\
 Data File : VX021821.D
 Acq On : 22 Apr 2021 16:56
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 IBLK

Quant Time: Apr 23 07:44:23 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042021W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 20 13:46:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	154696	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.830	114	281245	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.092	117	303834	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.061	152	163198	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.031	65	89307	50.89	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 101.78%			
35) Dibromofluoromethane	5.465	113	90644	50.40	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.80%			
50) Toluene-d8	8.696	98	348185	50.35	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 100.70%			
62) 4-Bromofluorobenzene	11.116	95	154574	59.85	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 119.70%			
Target Compounds						Qvalue
11) Tert butyl alcohol	3.014	59	1867	4.86	ug/l #	93
16) Acetone	2.428	43	1496	2.48	ug/l	94
17) Carbon Disulfide	2.556	76	1118	0.27	ug/l #	93

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

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