

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091521\
 Data File : VX024261.D
 Acq On : 15 Sep 2021 15:26
 Operator : JC/MD
 Sample : M3724-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 33127

Quant Time: Sep 16 04:44:25 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 14 12:18:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	128997	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	221985	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	210981	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94278	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.970	65	96213	52.305	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 104.620%			
35) Dibromofluoromethane	5.397	113	74514	50.052	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.100%			
50) Toluene-d8	8.653	98	271172	50.001	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 100.000%			
62) 4-Bromofluorobenzene	11.085	95	105922	50.186	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 100.380%			
Target Compounds						
16) Acetone	2.386	43	7506	9.570	ug/l	94
22) Diisopropyl ether	3.776	45	14000	3.298	ug/l #	87
29) Tetrahydrofuran	5.032	42	9916	13.714	ug/l #	71
31) Cyclohexane	5.483	56	10114	4.701	ug/l #	93
39) Methylcyclohexane	7.385	83	13623	5.591	ug/l	93
69) o-Xylene	10.646	106	71675	24.700	ug/l	97
80) 1,3,5-Trimethylbenzene	11.457	105	97749	16.518	ug/l	99
86) p-Isopropyltoluene	12.012	119	8533	1.420	ug/l	92

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

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