

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031323\
 Data File : VN076963.D
 Acq On : 13 Mar 2023 15:44
 Operator : JC\MD
 Sample : 01841-01DL2 400X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MORRIS-030823DL2

Quant Time: Mar 13 16:14:46 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031023W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 10 13:47:39 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.231	168	436535	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.107	114	701520	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.866	117	610839	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.795	152	267197	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.589	65	286545	43.987	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	87.980%		
35) Dibromofluoromethane	8.172	113	218733	46.784	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.560%		
50) Toluene-d8	10.572	98	820533	47.619	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	95.240%		
62) 4-Bromofluorobenzene	12.854	95	307017	48.713	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	97.420%		
Target Compounds						
8) Diethyl Ether	3.984	74	7301	2.214	ug/l	89
11) Tert butyl alcohol	5.537	59	16125	17.794	ug/l #	83
16) Acetone	4.449	43	16121	7.142	ug/l	93
40) Benzene	8.613	78	808169	37.916	ug/l	99
52) Toluene	10.630	92	536898	40.928	ug/l	99
67) Ethyl Benzene	11.966	91	55563	2.212	ug/l	96
68) m/p-Xylenes	12.072	106	130015	13.768	ug/l	86
69) o-Xylene	12.401	106	52724	5.751	ug/l	86
84) 1,2,4-Trimethylbenzene	13.483	105	36078	1.778	ug/l	91

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

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