

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022621\
 Data File : VN066056.D
 Acq On : 26 Feb 2021 17:06
 Operator : JC/MD
 Sample : M1488-05 20X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 41698

Quant Time: Feb 26 23:35:24 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 17 15:48:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.627	168	110162	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.550	114	177907	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.376	117	160127	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.315	152	64047	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.991	65	74766	48.53	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	97.06%		
35) Dibromofluoromethane	7.553	113	54770	49.74	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	99.48%		
50) Toluene-d8	10.058	98	215840	47.76	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	95.52%		
62) 4-Bromofluorobenzene	12.373	95	78121	46.04	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	92.08%		
Target Compounds						Qvalue
11) Tert butyl alcohol	4.750	59	6139	28.66	ug/l #	83
16) Acetone	3.788	43	4249462	8836.28	ug/l	96
20) Methylene Chloride	4.521	84	2235	1.56	ug/l #	90
25) 2-Butanone	6.791	43	42758	56.07	ug/l	98
29) Tetrahydrofuran	7.171	42	108531	206.73	ug/l	94
51) 4-Methyl-2-Pentanone	9.949	43	4795	3.12	ug/l	98
52) Toluene	10.122	92	80815	24.45	ug/l	95
67) Ethyl Benzene	11.482	91	5680	0.94	ug/l	96
68) m/p-Xylenes	11.588	106	9009	3.98	ug/l	91
69) o-Xylene	11.920	106	3699	1.70	ug/l	83

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

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