

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090222\
 Data File : VN074297.D
 Acq On : 02 Sep 2022 20:10
 Operator : JC\MD
 Sample : N4376-01DL 4X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-29DL

Quant Time: Sep 03 04:41:27 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090222W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 03 04:35:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	254636	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	450905	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	440837	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	203690	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	178621	45.000	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	90.000%		
35) Dibromofluoromethane	7.957	113	156281	52.304	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	104.600%		
50) Toluene-d8	10.380	98	547058	46.432	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	92.860%		
62) 4-Bromofluorobenzene	12.669	95	199756	46.542	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	93.080%		
Target Compounds						Qvalue
11) Tert butyl alcohol	5.298	59	20384	28.306	ug/l	95
16) Acetone	4.240	43	8049	4.852	ug/l #	88
19) Methyl tert-butyl Ether	5.534	73	38844	3.910	ug/l #	89
25) 2-Butanone	7.269	43	13128	4.686	ug/l	91
40) Benzene	8.392	78	237504	17.753	ug/l	94
51) 4-Methyl-2-Pentanone	10.275	43	19154	3.397	ug/l	91
52) Toluene	10.445	92	20354	2.462	ug/l	94
67) Ethyl Benzene	11.786	91	71940	4.050	ug/l	98
68) m/p-Xylenes	11.898	106	35511	5.355	ug/l	100
69) o-Xylene	12.222	106	13872	2.096	ug/l	98
84) 1,2,4-Trimethylbenzene	13.304	105	12015	0.829	ug/l	100
95) Naphthalene	15.439	128	8443	1.327	ug/l	95

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

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