

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040422\
 Data File : VN071838.D
 Acq On : 04 Apr 2022 17:53
 Operator : JC\MD
 Sample : N2200-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-29

Quant Time: Apr 05 01:20:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	613692	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1037047	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1077794	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	391430	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	343575	49.040	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	98.080%
35) Dibromofluoromethane	8.022	113	296064	48.516	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	97.040%
50) Toluene-d8	10.439	98	1314369	52.651	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	105.300%
62) 4-Bromofluorobenzene	12.727	95	442801	54.845	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	109.700%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	5802	1.938	ug/l	95
27) cis-1,2-Dichloroethene	7.328	96	4994	0.732	ug/l	82
30) Chloroform	7.810	83	7812	0.704	ug/l	97
38) Carbon Tetrachloride	8.204	117	7349	0.880	ug/l	92
44) Trichloroethene	9.216	130	5893	0.850	ug/l	97
64) Tetrachloroethene	10.975	164	45146	6.343	ug/l	94

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

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