

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042122\
 Data File : VX028249.D
 Acq On : 21 Apr 2022 18:19
 Operator : JC/MD
 Sample : N2459-20
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 4922B

Quant Time: Apr 22 06:17:19 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041922W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 19 13:38:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	413309	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	702310	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	659544	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	321935	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	289132	53.652	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 107.300%			
35) Dibromofluoromethane	5.391	113	267584	54.329	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 108.660%			
50) Toluene-d8	8.653	98	968162	51.180	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 102.360%			
62) 4-Bromofluorobenzene	11.079	95	351601	48.610	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 97.220%			
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	23136	12.731	ug/l	98
18) Methyl Acetate	2.709	43	3749	0.782	ug/l	98
20) Methylene Chloride	2.794	84	2354	0.589	ug/l #	76
25) 2-Butanone	4.562	43	19173	6.007	ug/l	91
30) Chloroform	5.105	83	8383	0.974	ug/l	97
51) 4-Methyl-2-Pentanone	8.580	43	12943	1.901	ug/l	96
52) Toluene	8.720	92	3291	0.252	ug/l	99
59) 2-Hexanone	9.439	43	2288	0.434	ug/l	93
64) Tetrachloroethene	9.275	164	1886	0.366	ug/l #	86
84) 1,2,4-Trimethylbenzene	11.756	105	4547	0.244	ug/l	97
95) Naphthalene	13.780	128	8427	0.395	ug/l #	93

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

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