

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042122\
 Data File : VX028248.D
 Acq On : 21 Apr 2022 17:56
 Operator : JC/MD
 Sample : N2459-18
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 4922A

Quant Time: Apr 22 06:16:51 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	383070	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	658416	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	635903	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	330255	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	279419	55.942	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 111.880%			
35) Dibromofluoromethane	5.391	113	252717	54.731	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 109.460%			
50) Toluene-d8	8.653	98	928595	52.361	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 104.720%			
62) 4-Bromofluorobenzene	11.079	95	345901	51.011	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 102.020%			
Target Compounds					Qvalue	
16) Acetone	2.373	43	433430	284.302	ug/l	93
18) Methyl Acetate	2.709	43	8500	1.914	ug/l	96
25) 2-Butanone	4.562	43	36144	12.219	ug/l	98
30) Chloroform	5.105	83	5304	0.665	ug/l	92
37) Ethyl Acetate	4.721	43	3178	0.520	ug/l #	87
40) Benzene	6.050	78	4298	0.244	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	22535	3.531	ug/l	97
52) Toluene	8.720	92	17533	1.431	ug/l	98
67) Ethyl Benzene	10.195	91	5208	0.229	ug/l	93
68) m/p-Xylenes	10.305	106	8194	0.886	ug/l	94
69) o-Xylene	10.646	106	4035	0.437	ug/l	93
84) 1,2,4-Trimethylbenzene	11.756	105	13217	0.690	ug/l	97
86) p-Isopropyltoluene	12.012	119	43856	2.142	ug/l	93
95) Naphthalene	13.780	128	72874	3.333	ug/l	98

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

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