

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010422\
 Data File : VX026270.D
 Acq On : 04 Jan 2022 14:41
 Operator : JC/MD
 Sample : M5335-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 20966

Quant Time: Jan 05 06:09:27 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	151689	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	263523	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	229325	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	100166	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	118762	54.118	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 108.240%		
35) Dibromofluoromethane	5.391	113	87351	50.986	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.980%		
50) Toluene-d8	8.653	98	308427	48.399	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 96.800%		
62) 4-Bromofluorobenzene	11.079	95	116388	51.304	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 102.600%		
Target Compounds						Qvalue
11) Tert butyl alcohol	2.952	59	9990	22.792	ug/l	97
16) Acetone	2.379	43	43471	36.561	ug/l	94
40) Benzene	6.043	78	87240	11.062	ug/l	98
51) 4-Methyl-2-Pentanone	8.579	43	8149	2.508	ug/l	99
52) Toluene	8.720	92	141130	29.068	ug/l	99
67) Ethyl Benzene	10.195	91	17567	1.950	ug/l	99
68) m/p-Xylenes	10.305	106	27218	8.093	ug/l	89
69) o-Xylene	10.640	106	9240	2.808	ug/l	88

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

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