

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX040422\
 Data File : VX027971.D
 Acq On : 04 Apr 2022 20:27
 Operator : JC/MD
 Sample : N2253-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 2822S

Quant Time: Apr 05 05:54:34 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032922W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 30 04:39:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	92425	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	149267	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	127374	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	58996	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	55671	46.683	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.360%
35) Dibromofluoromethane	5.391	113	47413	49.210	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.420%
50) Toluene-d8	8.653	98	171390	50.558	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.120%
62) 4-Bromofluorobenzene	11.079	95	68915	47.837	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	95.680%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.953	59	16016	69.917	ug/l	98
16) Acetone	2.373	43	445108	843.369	ug/l	90
17) Carbon Disulfide	2.514	76	3678	2.264	ug/l #	92
18) Methyl Acetate	2.703	43	7365	5.661	ug/l	90
20) Methylene Chloride	2.788	84	1222	1.287	ug/l	93
25) 2-Butanone	4.556	43	99220	129.201	ug/l	93
40) Benzene	6.044	78	1297	0.353	ug/l #	83
51) 4-Methyl-2-Pentanone	8.574	43	32844	19.649	ug/l	90
52) Toluene	8.720	92	2045	0.883	ug/l	99
59) 2-Hexanone	9.433	43	23124	17.967	ug/l	85
68) m/p-Xylenes	10.299	106	7258	4.335	ug/l	84
84) 1,2,4-Trimethylbenzene	11.756	105	1136	0.299	ug/l	78

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

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