

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032822\
 Data File : VX027756.D
 Acq On : 28 Mar 2022 17:42
 Operator : JC/MD
 Sample : N2154-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-4

Quant Time: Mar 29 05:40:24 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032522W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 25 16:26:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	68362	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	121410	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	110337	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	48163	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	50382	44.935	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	89.860%
35) Dibromofluoromethane	5.391	113	39875	50.013	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.020%
50) Toluene-d8	8.653	98	148546	48.197	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.400%
62) 4-Bromofluorobenzene	11.079	95	59095	47.253	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	94.500%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.953	59	47478	227.427	ug/l	99
16) Acetone	2.379	43	26084	55.007	ug/l	92
17) Carbon Disulfide	2.514	76	1350	0.693	ug/l	# 77
25) 2-Butanone	4.562	43	6487	8.518	ug/l	92
30) Chloroform	5.099	83	2225	1.432	ug/l	# 81
45) 1,2-Dichloropropane	7.440	63	2359	2.584	ug/l	99
51) 4-Methyl-2-Pentanone	8.580	43	4399	2.895	ug/l	89
68) m/p-Xylenes	10.311	106	866	0.552	ug/l	86
69) o-Xylene	10.646	106	477	0.302	ug/l	86

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

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