

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060622\
 Data File : VX029235.D
 Acq On : 06 Jun 2022 13:37
 Operator : JC/MD
 Sample : N3174-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MID0622

Quant Time: Jun 07 06:04:01 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 01 04:45:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	317019	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	592378	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	580436	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	270915	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	214993	60.141	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 120.280%	
35) Dibromofluoromethane	5.391	113	197996	50.320	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 100.640%	
50) Toluene-d8	8.653	98	705245	49.006	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 98.020%	
62) 4-Bromofluorobenzene	11.085	95	283669	50.994	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 101.980%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.965	59	16160	19.514	ug/l	97
16) Acetone	2.386	43	4808	3.416	ug/l	91
17) Carbon Disulfide	2.520	76	3903	0.498	ug/l	99
19) Methyl tert-butyl Ether	3.123	73	6786	0.700	ug/l #	85
31) Cyclohexane	5.476	56	12288	2.644	ug/l	93
40) Benzene	6.037	78	81949	5.465	ug/l	97

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

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