

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060222\
 Data File : VX029139.D
 Acq On : 02 Jun 2022 16:42
 Operator : JC/MD
 Sample : N3022-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 22BR-20220525

Quant Time: Jun 03 04:32:44 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	271781	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	498596	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	476872	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	221086	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	177024	57.762	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	115.520%
35) Dibromofluoromethane	5.391	113	163617	49.404	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.800%
50) Toluene-d8	8.653	98	590777	48.774	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.540%
62) 4-Bromofluorobenzene	11.085	95	233629	49.898	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	99.800%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	3158	2.617	ug/l #	85
20) Methylene Chloride	2.788	84	1469	0.522	ug/l #	95
27) cis-1,2-Dichloroethene	4.495	96	13290	4.033	ug/l	89
30) Chloroform	5.105	83	1990	0.394	ug/l	86
44) Trichloroethene	7.129	130	6701	1.104	ug/l	92

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

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