

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060322\
 Data File : VX029189.D
 Acq On : 03 Jun 2022 14:52
 Operator : JC/MD
 Sample : N3164-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 27BR-20220531

Quant Time: Jun 06 02:35:09 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	308357	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	571059	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	551945	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	254379	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	201079	57.828	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	115.660%
35) Dibromofluoromethane	5.391	113	190424	50.203	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.400%
50) Toluene-d8	8.653	98	673747	48.565	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.140%
62) 4-Bromofluorobenzene	11.085	95	265835	49.572	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	99.140%
Target Compounds						
						Qvalue
3) Chloromethane	1.295	50	649	0.231	ug/l	97
16) Acetone	2.380	43	5674	4.144	ug/l #	85
23) Vinyl Acetate	3.733	43	2543	0.365	ug/l	100
29) Tetrahydrofuran	5.026	42	2728	1.926	ug/l	91
37) Ethyl Acetate	4.715	43	575228	121.612	ug/l	99

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

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