

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060322\
 Data File : VX029191.D
 Acq On : 03 Jun 2022 15:39
 Operator : JC/MD
 Sample : N3164-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 31BR-20220531

Quant Time: Jun 06 02:35:28 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	310582	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	582009	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	562227	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	262002	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	203506	58.107	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	116.220%
35) Dibromofluoromethane	5.385	113	191662	49.578	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.160%
50) Toluene-d8	8.653	98	688090	48.666	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.340%
62) 4-Bromofluorobenzene	11.079	95	273715	50.081	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	100.160%
Target Compounds						
						Qvalue
3) Chloromethane	1.289	50	914	0.323	ug/l	90
16) Acetone	2.386	43	12052	8.739	ug/l	97
24) 1,1-Dichloroethane	3.611	63	4451	0.856	ug/l	96
27) cis-1,2-Dichloroethene	4.489	96	15693	4.167	ug/l	86
29) Tetrahydrofuran	5.013	42	1994	1.398	ug/l #	83
37) Ethyl Acetate	4.721	43	93411	19.377	ug/l	98
44) Trichloroethene	7.129	130	21929	4.380	ug/l	98

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

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