

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060322\
 Data File : VX029192.D
 Acq On : 03 Jun 2022 16:02
 Operator : JC/MD
 Sample : N3164-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 31S-20220531

Quant Time: Jun 06 02:35:37 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	303028	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	571339	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	551988	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	255464	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	200805	58.765	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 117.540%		
35) Dibromofluoromethane	5.385	113	189484	49.930	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 99.860%		
50) Toluene-d8	8.653	98	675229	48.648	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 97.300%		
62) 4-Bromofluorobenzene	11.079	95	268287	50.004	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 100.000%		
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	3989	2.965	ug/l	# 84
24) 1,1-Dichloroethane	3.611	63	5775	1.138	ug/l	# 95
27) cis-1,2-Dichloroethene	4.501	96	11655	3.172	ug/l	87
29) Tetrahydrofuran	5.032	42	2340	1.681	ug/l	# 71
37) Ethyl Acetate	4.715	43	111781	23.621	ug/l	99
44) Trichloroethene	7.141	130	4692	0.398	ug/l	93

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

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