

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032222\
 Data File : VX027627.D
 Acq On : 22 Mar 2022 20:13
 Operator : JC/MD
 Sample : N2028-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-118

Quant Time: Mar 23 01:51:14 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 17 15:02:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	354647	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	572129	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	504378	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	231037	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	189607	43.399	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	86.800%		
35) Dibromofluoromethane	5.391	113	177831	48.311	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.620%		
50) Toluene-d8	8.653	98	659507	47.347	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	94.700%		
62) 4-Bromofluorobenzene	11.085	95	229162	44.760	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	89.520%		
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.965	59	4870	5.849	ug/l	96
16) Acetone	2.386	43	13510	7.587	ug/l	100
17) Carbon Disulfide	2.514	76	8397	1.002	ug/l	97
40) Benzene	6.037	78	9653	0.669	ug/l #	88
51) 4-Methyl-2-Pentanone	8.580	43	4704	0.913	ug/l	96
52) Toluene	8.726	92	5638	0.583	ug/l	99
67) Ethyl Benzene	10.195	91	3893	0.219	ug/l #	89
68) m/p-Xylenes	10.305	106	6251	0.894	ug/l	96
69) o-Xylene	10.646	106	1787	0.263	ug/l	94
84) 1,2,4-Trimethylbenzene	11.756	105	5425	0.384	ug/l	94

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

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