

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101122\
 Data File : VX032062.D
 Acq On : 12 Oct 2022 05:28
 Operator : JC/MD
 Sample : N5067-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B2(19-20)

Quant Time: Oct 12 06:23:59 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 06 14:28:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	77975	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	118052	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	128365	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	73026	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	82523	51.245	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.500%			
35) Dibromofluoromethane	5.391	113	62917	47.084	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 94.160%			
50) Toluene-d8	8.647	98	242005	50.330	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 100.660%			
62) 4-Bromofluorobenzene	11.079	95	80771	47.515	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 95.040%			
Target Compounds					Qvalue	
16) Acetone	2.386	43	28408	51.906	ug/l	99
20) Methylene Chloride	2.794	84	9656	6.714	ug/l	88
25) 2-Butanone	4.574	43	5374	6.281	ug/l	95
37) Ethyl Acetate	4.715	43	11888	7.126	ug/l	98
40) Benzene	6.044	78	31717	6.113	ug/l	98
43) Isopropyl Acetate	6.342	43	56790	22.437	ug/l	99
67) Ethyl Benzene	10.195	91	25391	3.859	ug/l	97
68) m/p-Xylenes	10.305	106	4344	1.661	ug/l	94
73) Isopropylbenzene	10.963	105	7683	1.395	ug/l	94
84) 1,2,4-Trimethylbenzene	11.756	105	10928	2.355	ug/l	98
90) Hexachloroethane	12.701	117	2008	3.918	ug/l #	15
95) Naphthalene	13.774	128	165016	30.893	ug/l	99

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

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