

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
 Data File : VN072767.D
 Acq On : 10 Jun 2022 16:10
 Operator : JC\MD
 Sample : N3293-16 50X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 40809

Quant Time: Jun 11 01:54:56 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 09 05:17:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	167100	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.963	114	317001	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	274703	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	101354	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	253141	56.444	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 112.880%	
35) Dibromofluoromethane	8.016	113	183914	58.633	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 117.260%	
50) Toluene-d8	10.439	98	573169	45.800	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 91.600%	
62) 4-Bromofluorobenzene	12.727	95	209201	48.715	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 97.440%	
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	70151	32.643	ug/l	88
40) Benzene	8.457	78	6912	0.523	ug/l	98
52) Toluene	10.504	92	33065	4.008	ug/l	98
67) Ethyl Benzene	11.839	91	14436	1.049	ug/l	# 87
68) m/p-Xylenes	11.945	106	19386	3.963	ug/l	77
69) o-Xylene	12.280	106	14051	2.858	ug/l	88
78) n-propylbenzene	12.916	91	3436	0.289	ug/l	88
80) 1,3,5-Trimethylbenzene	13.057	105	6012	0.698	ug/l	96
84) 1,2,4-Trimethylbenzene	13.363	105	30859	3.744	ug/l	93
95) Naphthalene	15.504	128	30492	4.579	ug/l	99

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

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