

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060222\
 Data File : VU048804.D
 Acq On : 02 Jun 2022 15:25
 Operator : SY/MD
 Sample : N3087-06DL 20X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-06DL

Quant Time: Jun 02 15:54:24 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 27 06:42:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.369	168	351642	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.247	114	585590	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	528696	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	283044	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.697	65	235798	44.650	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	89.300%		
35) Dibromofluoromethane	5.285	113	185022	44.288	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	88.580%		
50) Toluene-d8	7.896	98	705746	47.126	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	94.260%		
62) 4-Bromofluorobenzene	10.632	95	273020	48.206	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	96.420%		
Target Compounds					Qvalue	
16) Acetone	2.626	43	10071	3.742	ug/l	99
25) 2-Butanone	4.710	43	15072	3.898	ug/l	100
39) Methylcyclohexane	6.758	83	17808	2.785	ug/l	98
40) Benzene	5.771	78	749750	41.650	ug/l	100
52) Toluene	7.970	92	66010	6.131	ug/l	100
67) Ethyl Benzene	9.568	91	400447	21.233	ug/l	99
68) m/p-Xylenes	9.693	106	176152	23.881	ug/l	98
69) o-Xylene	10.099	106	44971	6.206	ug/l	98
73) Isopropylbenzene	10.484	105	24757	1.355	ug/l	98
78) n-propylbenzene	10.906	91	48037	2.206	ug/l	98
80) 1,3,5-Trimethylbenzene	11.086	105	54309	3.418	ug/l	99
84) 1,2,4-Trimethylbenzene	11.468	105	369594	22.255	ug/l	99

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

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