

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060222\
 Data File : VU048805.D
 Acq On : 02 Jun 2022 15:48
 Operator : SY/MD
 Sample : N3087-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-04

Quant Time: Jun 02 16:46:32 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	5.372	168	334050	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.247	114	557746	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	508353	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	259480	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	223120	44.475	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	88.940%
35) Dibromofluoromethane	5.285	113	176017	44.236	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	88.480%
50) Toluene-d8	7.896	98	669279	46.922	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	93.840%
62) 4-Bromofluorobenzene	10.632	95	250514	46.440	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.880%
Target Compounds						
					Qvalue	
16) Acetone	2.626	43	7520	2.941	ug/l	93
25) 2-Butanone	4.716	43	6493	1.768	ug/l	97
30) Chloroform	5.086	83	1762	0.231	ug/l	85
39) Methylcyclohexane	6.761	83	1673	0.275	ug/l #	73
40) Benzene	5.774	78	27846	1.624	ug/l	98
52) Toluene	7.967	92	23646	2.306	ug/l	98
67) Ethyl Benzene	9.571	91	31444	1.734	ug/l	98
68) m/p-Xylenes	9.693	106	54314	7.658	ug/l	98
69) o-Xylene	10.099	106	20797	2.985	ug/l	97
80) 1,3,5-Trimethylbenzene	11.089	105	7727	0.530	ug/l	99
84) 1,2,4-Trimethylbenzene	11.468	105	26065	2.808	ug/l	96
95) Naphthalene	14.092	128	21185	3.517	ug/l	97

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

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