

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010623\
 Data File : VN076142.D
 Acq On : 06 Jan 2023 18:36
 Operator : JC\MD
 Sample : 01066-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OILY-WATER-1

Quant Time: Jan 06 22:57:52 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 14 13:03:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	336635	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	537499	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	462098	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	191866	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	177348	48.426	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.860%
35) Dibromofluoromethane	8.177	113	161697	48.577	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.160%
50) Toluene-d8	10.571	98	616805	49.561	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.120%
62) 4-Bromofluorobenzene	12.853	95	189849	46.057	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.120%
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	2826	2.238	ug/l	91
52) Toluene	10.635	92	4512	0.508	ug/l	93
68) m/p-Xylenes	12.071	106	4506	0.711	ug/l	86
69) o-Xylene	12.394	106	4754	0.758	ug/l	93
80) 1,3,5-Trimethylbenzene	13.182	105	5797	0.500	ug/l	85
84) 1,2,4-Trimethylbenzene	13.482	105	10634	0.924	ug/l	93

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

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