

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052423\
 Data File : VN077860.D
 Acq On : 24 May 2023 18:19
 Operator : JC\MD
 Sample : 02897-02 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OILY-WATER-0523

Quant Time: May 25 01:43:54 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051523W.M
 Quant Title : SW846 8260
 QLast Update : Tue May 16 04:07:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	414342	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	737850	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.876	117	672686	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	256845	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	268485	43.307	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.620%
35) Dibromofluoromethane	8.171	113	230793	47.968	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.940%
50) Toluene-d8	10.576	98	910629	49.171	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.340%
62) 4-Bromofluorobenzene	12.859	95	323328	49.433	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.860%
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	54493	27.938	ug/l	89
18) Methyl Acetate	5.053	43	12400	1.515	ug/l #	71
27) cis-1,2-Dichloroethene	7.500	96	10193	1.737	ug/l	99
29) Tetrahydrofuran	7.859	42	11059	5.535	ug/l #	81
37) Ethyl Acetate	7.571	43	50200	8.003	ug/l	96
44) Trichloroethene	9.359	130	52671	10.016	ug/l	94
64) Tetrachloroethene	11.112	164	1688	0.359	ug/l #	76
68) m/p-Xylenes	12.070	106	7486	0.756	ug/l	97
69) o-Xylene	12.400	106	3155	0.324	ug/l	85
80) 1,3,5-Trimethylbenzene	13.182	105	8562	0.460	ug/l	96
84) 1,2,4-Trimethylbenzene	13.488	105	26482	1.447	ug/l	92
86) p-Isopropyltoluene	13.735	119	1938037	105.298	ug/l	99

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

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