

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111522\
 Data File : VN075283.D
 Acq On : 15 Nov 2022 14:57
 Operator : JC\MD
 Sample : N5598-13
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FRAC-TANK-SLURRY-0760533

Quant Time: Nov 16 01:03:00 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 09:48:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	158758	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	289210	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	266346	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	111268	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	106429	49.755	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.500%
35) Dibromofluoromethane	8.171	113	87758	51.689	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.380%
50) Toluene-d8	10.565	98	346791	52.647	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	105.300%
62) 4-Bromofluorobenzene	12.847	95	114222	49.363	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.720%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	10624	11.730	ug/l	98
20) Methylene Chloride	5.283	84	560274	277.435	ug/l	95
37) Ethyl Acetate	7.565	43	17051	5.915	ug/l	98
43) Isopropyl Acetate	8.688	43	72937	16.312	ug/l #	92

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

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