

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091922\
 Data File : VN074582.D
 Acq On : 19 Sep 2022 13:36
 Operator : JC\MD
 Sample : N4647-06DL 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1037-MW-11(15)DL

Quant Time: Sep 20 04:19:15 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091522W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 19:20:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	282029	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	537197	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	526101	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	227041	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	220257	46.028	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	92.060%
35) Dibromofluoromethane	7.957	113	180386	50.549	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.100%
50) Toluene-d8	10.381	98	657779	47.663	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.320%
62) 4-Bromofluorobenzene	12.669	95	245567	51.953	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.900%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	6358	2.076	ug/l #	79
27) cis-1,2-Dichloroethene	7.245	96	20117	3.852	ug/l	94
44) Trichloroethene	9.151	130	47661	11.277	ug/l	87
64) Tetrachloroethene	10.916	164	109266	30.263	ug/l	91

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

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