

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082447.D
 Acq On : 03 Jun 2024 20:29
 Operator : JC\MD
 Sample : P2686-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPQ

Quant Time: Jun 04 00:35:50 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	153045	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	285810	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	285714	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	117076	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	135240	55.962	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.920%
35) Dibromofluoromethane	8.171	113	97174	55.270	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	110.540%
50) Toluene-d8	10.571	98	379333	56.275	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	112.560%
62) 4-Bromofluorobenzene	12.853	95	128240	53.302	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	=	106.600%
Target Compounds						
						Qvalue
16) Acetone	4.436	43	91988	92.792	ug/l	94
20) Methylene Chloride	5.283	84	13777	6.888	ug/l #	73
25) 2-Butanone	7.494	43	14414	9.590	ug/l #	79
43) Isopropyl Acetate	8.694	43	275293	47.080	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	16969	5.108	ug/l #	72

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

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