

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082448.D
 Acq On : 03 Jun 2024 20:53
 Operator : JC\MD
 Sample : P2686-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPP

Quant Time: Jun 04 00:36:17 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	8.229	168	155237	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	291411	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	292223	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	120867	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	136233	55.577	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 111.160%	
35) Dibromofluoromethane	8.171	113	98749	55.086	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 110.180%	
50) Toluene-d8	10.570	98	386378	56.218	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 112.440%	
62) 4-Bromofluorobenzene	12.853	95	129487	52.786	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 105.580%	
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	104677	104.100	ug/l	90
20) Methylene Chloride	5.288	84	13714	6.760	ug/l #	75
25) 2-Butanone	7.488	43	18873	12.379	ug/l #	80
40) Benzene	8.612	78	59460	6.632	ug/l	97
43) Isopropyl Acetate	8.694	43	270230	45.326	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	18093	5.341	ug/l #	73

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

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