

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060524\
 Data File : VN082470.D
 Acq On : 05 Jun 2024 15:01
 Operator : JC\MD
 Sample : P2704-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPO

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/06/2024
 Supervised By : Mahesh Dadoda 06/06/2024

Quant Time: Jun 06 00:57:16 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	8.224	168	159427	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	299034	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	304559	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	125329	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	141443	56.186	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 112.380%			
35) Dibromofluoromethane	8.171	113	101797	55.339	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 110.680%			
50) Toluene-d8	10.570	98	401541	56.935	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 113.880%#			
62) 4-Bromofluorobenzene	12.853	95	130013	51.649	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 103.300%			
Target Compounds						
16) Acetone	4.430	43	97803	94.708	ug/l	93
20) Methylene Chloride	5.277	84	4475	2.148	ug/l #	78
25) 2-Butanone	7.494	43	23973	15.311	ug/l #	77
43) Isopropyl Acetate	8.688	43	294177	48.085	ug/l #	70
51) 4-Methyl-2-Pentanone	10.453	43	22792m	6.557	ug/l	

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

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