

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081924\
 Data File : VN083375.D
 Acq On : 19 Aug 2024 16:16
 Operator : JC\MD
 Sample : P3629-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-4

Quant Time: Aug 20 04:47:28 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:30:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	123808	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	252966	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	274805	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	119589	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	108907	61.799	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	123.600%
35) Dibromofluoromethane	8.165	113	87615	55.489	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	110.980%
50) Toluene-d8	10.565	98	330019	56.031	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	112.060%
62) 4-Bromofluorobenzene	12.847	95	135420	58.975	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	117.960%
Target Compounds						
						Qvalue
16) Acetone	4.442	43	32600	47.276	ug/l	96
20) Methylene Chloride	5.277	84	7730	4.871	ug/l #	72
25) 2-Butanone	7.500	43	6097	5.758	ug/l	95
43) Isopropyl Acetate	8.688	43	125203	25.777	ug/l #	92
51) 4-Methyl-2-Pentanone	10.447	43	94910	37.529	ug/l	97

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

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