

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020724\
 Data File : VN080941.D
 Acq On : 07 Feb 2024 17:15
 Operator : JC\MD
 Sample : PB158850ZHE#05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB158850ZHE#05

Quant Time: Feb 08 00:28:40 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020624W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 07 06:22:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	271592	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	592188	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	752703	50.000	ug/l	# 0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	318753	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	193826	56.573	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	113.140%
35) Dibromofluoromethane	8.165	113	180316	56.745	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	113.480%
50) Toluene-d8	10.571	98	721175	62.744	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	125.480%#
62) 4-Bromofluorobenzene	12.853	95	311309	74.557	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	149.120%#
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	45461	40.331	ug/l #	83
20) Methylene Chloride	5.271	84	8322	2.437	ug/l #	74
37) Ethyl Acetate	7.565	43	7982	2.231	ug/l #	88
43) Isopropyl Acetate	8.694	43	294585	44.786	ug/l #	90
51) 4-Methyl-2-Pentanone	10.453	43	21179	5.343	ug/l #	78

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

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