

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072723\
 Data File : VN078666.D
 Acq On : 27 Jul 2023 17:44
 Operator : JC\MD
 Sample : PB154443ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154443ZHE#03

Quant Time: Jul 28 01:10:00 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	409809	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	9.112	114	733657	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	695756	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	255843	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	268225	48.822	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	97.640%	
35) Dibromofluoromethane	8.177	113	263294	52.867	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	105.740%	
50) Toluene-d8	10.576	98	908432	48.285	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	96.560%	
62) 4-Bromofluorobenzene	12.853	95	304074	52.471	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	104.940%	
Target Compounds						
						Qvalue
16) Acetone	4.453	43	94929	55.470	ug/l	97
20) Methylene Chloride	5.289	84	26684	4.319	ug/l	98
25) 2-Butanone	7.500	43	50192	18.632	ug/l #	83
37) Ethyl Acetate	7.588	43	18049	3.011	ug/l #	90
43) Isopropyl Acetate	8.700	43	435356	44.112	ug/l #	77
51) 4-Methyl-2-Pentanone	10.453	43	175243	29.852	ug/l	89

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

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