

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081122\
 Data File : VN073827.D
 Acq On : 11 Aug 2022 16:03
 Operator : JC\MD
 Sample : PB146912ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB146912ZHE#02

Quant Time: Aug 12 06:03:28 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 10 05:09:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.022	168	381578	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	615795	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	618508	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	288789	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	226405	48.688	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.380%
35) Dibromofluoromethane	7.957	113	214991	55.599	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	111.200%
50) Toluene-d8	10.380	98	721602	47.423	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.840%
62) 4-Bromofluorobenzene	12.674	95	270976	50.689	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.380%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	81237	32.526	ug/l	95
18) Methyl Acetate	4.787	43	6559	1.056	ug/l	93
20) Methylene Chloride	5.016	84	16995	2.095	ug/l	96
25) 2-Butanone	7.269	43	25812	6.655	ug/l	96
37) Ethyl Acetate	7.345	43	70842	9.761	ug/l	99
43) Isopropyl Acetate	8.498	43	277971	28.355	ug/l #	92

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

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