

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072922\
 Data File : VN073643.D
 Acq On : 29 Jul 2022 18:37
 Operator : JC\MD
 Sample : PB146545ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB146545ZHE#04

Quant Time: Aug 01 04:08:09 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 29 02:20:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	340497	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	563391	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	553623	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	237887	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	191763	46.213	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	92.420%
35) Dibromofluoromethane	7.951	113	179578	50.761	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.520%
50) Toluene-d8	10.380	98	629762	45.237	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.480%
62) 4-Bromofluorobenzene	12.674	95	225858	48.789	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.580%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	104540	46.905	ug/l	98
18) Methyl Acetate	4.787	43	6921	1.248	ug/l #	88
20) Methylene Chloride	5.016	84	23781	4.594	ug/l	95
25) 2-Butanone	7.269	43	31377	9.065	ug/l	96
37) Ethyl Acetate	7.345	43	58968	8.880	ug/l	99
43) Isopropyl Acetate	8.498	43	238645	26.608	ug/l #	92

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

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