

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080222\
 Data File : VX030355.D
 Acq On : 02 Aug 2022 16:18
 Operator : JC/MD
 Sample : PB146634ZHE#02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB146634ZHE#02

Quant Time: Aug 03 01:08:49 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 25 16:49:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	157508	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	286681	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	284776	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	160939	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	164854	52.073	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 104.140%		
35) Dibromofluoromethane	5.397	113	133034	48.543	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 97.080%		
50) Toluene-d8	8.653	98	505237	47.605	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 95.220%		
62) 4-Bromofluorobenzene	11.085	95	179458	45.562	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 91.120%		
Target Compounds						
						Qvalue
16) Acetone	2.380	43	101197	62.436	ug/l	96
20) Methylene Chloride	2.794	84	16835	3.712	ug/l	97
25) 2-Butanone	4.568	43	32994	12.114	ug/l	95
29) Tetrahydrofuran	5.025	42	12103	6.604	ug/l	96
37) Ethyl Acetate	4.727	43	51416	9.206	ug/l	97
43) Isopropyl Acetate	6.348	43	218601	26.700	ug/l	97

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

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