

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071024\
 Data File : VX042267.D
 Acq On : 10 Jul 2024 15:10
 Operator : JC/MD
 Sample : PB161974ZHE#08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB161974ZHE#08

Quant Time: Jul 11 00:57:57 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 21 23:51:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	139022	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	223854	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	200756	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	85369	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	82341	51.364	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	102.720%
35) Dibromofluoromethane	5.385	113	78422	49.373	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.740%
50) Toluene-d8	8.647	98	244527	43.822	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	87.640%#
62) 4-Bromofluorobenzene	11.079	95	83866	40.773	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	81.540%#
Target Compounds						
						Qvalue
16) Acetone	2.392	43	25059	35.599	ug/l	93
18) Methyl Acetate	2.709	43	4649	2.421	ug/l	98
20) Methylene Chloride	2.788	84	8231	5.340	ug/l #	86
37) Ethyl Acetate	4.745	43	3106	1.414	ug/l #	82
43) Isopropyl Acetate	6.348	43	110104	32.647	ug/l #	93

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

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