

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036407.D
 Acq On : 05 Jul 2023 16:31
 Operator : JC/MD
 Sample : PB153983ZHE#08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983ZHE#08

Quant Time: Jul 06 04:53:38 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 23 15:03:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	167020	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	281906	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	249185	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	117018	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	107667	44.972	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	89.940%		
35) Dibromofluoromethane	5.385	113	89133	47.952	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.900%		
50) Toluene-d8	8.647	98	340914	49.104	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.200%		
62) 4-Bromofluorobenzene	11.079	95	120098	49.111	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	98.220%		
Target Compounds						
					Qvalue	
16) Acetone	2.398	43	220976	215.942	ug/l	96
18) Methyl Acetate	2.715	43	4669	1.110	ug/l	98
20) Methylene Chloride	2.788	84	7471	3.391	ug/l	97
25) 2-Butanone	4.574	43	117525	79.933	ug/l	97
37) Ethyl Acetate	4.727	43	55321	19.589	ug/l	98
43) Isopropyl Acetate	6.342	43	203063	45.074	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	533202	183.706	ug/l	87

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

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