

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
 Data File : VX043121.D
 Acq On : 24 Sep 2024 14:39
 Operator : JC/MD
 Sample : PB163627TB
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB163627TB

Quant Time: Sep 25 01:28:39 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 24 03:13:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	92238	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	180360	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	172332	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	91350	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	80583	52.596	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	105.200%
35) Dibromofluoromethane	5.385	113	62082	48.794	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.580%
50) Toluene-d8	8.647	98	224440	52.140	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	104.280%
62) 4-Bromofluorobenzene	11.079	95	93055	53.681	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	107.360%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	44426	62.003	ug/l	99
18) Methyl Acetate	2.703	43	176265	91.399	ug/l	99
20) Methylene Chloride	2.794	84	6960	6.106	ug/l	92
51) 4-Methyl-2-Pentanone	8.580	43	6993	3.470	ug/l	99

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

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