

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036486.D
 Acq On : 10 Jul 2023 16:31
 Operator : JC/MD
 Sample : PB154051ZHE#08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB154051ZHE#08

Quant Time: Jul 11 01:30:56 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	5.550	168	167801	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	289530	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	268199	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	124024	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	113731	47.283	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	94.560%	
35) Dibromofluoromethane	5.385	113	90809	47.567	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.140%	
50) Toluene-d8	8.647	98	359756	50.454	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	100.900%	
62) 4-Bromofluorobenzene	11.079	95	128228	51.055	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	102.100%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	84547	82.236	ug/l	93
20) Methylene Chloride	2.794	84	7352	3.321	ug/l	87
25) 2-Butanone	4.556	43	58504	39.605	ug/l	93
37) Ethyl Acetate	4.715	43	16165	5.573	ug/l	98
43) Isopropyl Acetate	6.342	43	150379	32.501	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	290642	97.499	ug/l #	81

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

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