

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD102023\
 Data File : VD077373.D
 Acq On : 20 Oct 2023 19:08
 Operator : JC/SY
 Sample : 04936-07RE
 Misc : 5.08G/5.0ml/MSVOA_D/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 1084RE

Quant Time: Oct 21 05:32:38 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D101723S.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 18 03:02:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.881	168	25549	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.781	114	58164	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.581	117	38353	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.516	152	7720	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.228	65	31867	139.683	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 279.360%#	
35) Dibromofluoromethane	7.810	113	29583	72.396	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 144.800%	
50) Toluene-d8	10.269	98	41039	28.043	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	= 56.080%#	
62) 4-Bromofluorobenzene	12.575	95	8446	20.049	ug/l	0.00
Spiked Amount	50.000	Range	30 - 143	Recovery	= 40.100%	
Target Compounds						
					Qvalue	
16) Acetone	4.016	43	2967350	59314.092	ug/l	97
20) Methylene Chloride	4.805	84	6440	18.239	ug/l #	87
52) Toluene	10.334	92	859139	807.556	ug/l	97
67) Ethyl Benzene	11.681	91	4417	3.187	ug/l #	85
68) m/p-Xylenes	11.787	106	2489	4.374	ug/l	100
69) o-Xylene	12.122	106	933	1.781	ug/l	88

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

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