

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041221\
 Data File : VU043031.D
 Acq On : 12 Apr 2021 18:19
 Operator : SY/MD
 Sample : M1946-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FS-SB10

Quant Time: Apr 13 05:15:18 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U041221W.M
 Quant Title : SW846 8260
 QLast Update : Mon Apr 12 13:17:29 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	117627	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	193495	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	188347	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	94291	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	98440	49.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.56%	
35) Dibromofluoromethane	5.298	113	69372	47.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.42%	
50) Toluene-d8	7.906	98	248498	48.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.60%	
62) 4-Bromofluorobenzene	10.639	95	91341	44.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.68%	
Target Compounds						
						Qvalue
16) Acetone	2.639	43	2121	1.52	ug/l	99
40) Benzene	5.780	78	7577	1.29	ug/l	91
52) Toluene	7.980	92	2017	0.55	ug/l	93
67) Ethyl Benzene	9.578	91	4119	0.57	ug/l	94
68) m/p-Xylenes	9.697	106	3448	1.28	ug/l	100
69) o-Xylene	10.105	106	1351	0.51	ug/l	81
84) 1,2,4-Trimethylbenzene	11.472	105	1484	0.26	ug/l	85
95) Naphthalene	14.095	128	7798	3.82	ug/l	97

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

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