

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092721\
 Data File : VU044898.D
 Acq On : 27 Sep 2021 18:52
 Operator : SY/MD
 Sample : M3888-11DL 100X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 40458DL

Quant Time: Sep 27 19:12:46 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092421W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 25 09:30:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	197185	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	350428	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	327286	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	165353	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	155020	45.045	ug/l	0.00
Spiked Amount	50.000		Recovery	=	90.100%	
35) Dibromofluoromethane	5.292	113	105889	44.047	ug/l	0.00
Spiked Amount	50.000		Recovery	=	88.100%	
50) Toluene-d8	7.903	98	439108	48.363	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.720%	
62) 4-Bromofluorobenzene	10.636	95	178881	53.045	ug/l	0.00
Spiked Amount	50.000		Recovery	=	106.080%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	26614	13.710	ug/l	97
40) Benzene	5.777	78	95938	8.463	ug/l	99
52) Toluene	7.973	92	1116773	164.780	ug/l	99
67) Ethyl Benzene	9.575	91	297287	24.514	ug/l	99
68) m/p-Xylenes	9.697	106	490544	105.761	ug/l	97
69) o-Xylene	10.105	106	226410	50.481	ug/l	99
73) Isopropylbenzene	10.488	105	12823	1.093	ug/l	99
78) n-propylbenzene	10.909	91	38188	2.685	ug/l	100
80) 1,3,5-Trimethylbenzene	11.092	105	77140	7.760	ug/l	99
84) 1,2,4-Trimethylbenzene	11.472	105	292066	29.801	ug/l	99
95) Naphthalene	14.089	128	61427	5.695	ug/l	100

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

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