

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040921\
 Data File : VU043001.D
 Acq On : 09 Apr 2021 14:46
 Operator : SY/MD
 Sample : M1977-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GWBR3-200-220-040721

Quant Time: Apr 09 15:44:35 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U040921W.M
 Quant Title : SW846 8260
 QLast Update : Fri Apr 09 09:26:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	120582	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	199876	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	193095	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	97855	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	99143	50.30	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.60%	
35) Dibromofluoromethane	5.298	113	69914	47.90	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.80%	
50) Toluene-d8	7.906	98	251473	47.83	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.66%	
62) 4-Bromofluorobenzene	10.639	95	94513	44.00	ug/l	0.00
Spiked Amount	50.000		Recovery	=	88.00%	
Target Compounds						
						Qvalue
12) 1,1-Dichloroethene	2.584	96	21555	19.41	ug/l	97
16) Acetone	2.626	43	10122	9.16	ug/l	93
20) Methylene Chloride	3.054	84	1121	0.79	ug/l	91
21) trans-1,2-Dichloroethene	3.359	96	1055	0.84	ug/l #	75
24) 1,1-Dichloroethane	3.880	63	5196	2.09	ug/l #	95
25) 2-Butanone	4.713	43	6807	3.85	ug/l	100
27) cis-1,2-Dichloroethene	4.674	96	14832	10.51	ug/l	90
30) Chloroform	5.096	83	8294	3.09	ug/l	100
42) 1,2-Dichloroethane	5.797	62	4278	1.68	ug/l	91
44) Trichloroethene	6.549	130	53229	35.90	ug/l	93
51) 4-Methyl-2-Pentanone	7.803	43	3779	10.45	ug/l	98
52) Toluene	7.976	92	466299	129.38	ug/l	100
59) 2-Hexanone	8.694	43	3474	8.94	ug/l	99
84) 1,2,4-Trimethylbenzene	11.481	105	2053	0.36	ug/l #	51
95) Naphthalene	14.095	128	5510	3.94	ug/l #	88

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

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