

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092222\
 Data File : VX031604.D
 Acq On : 23 Sep 2022 05:40
 Operator : JC/MD
 Sample : N4750-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SB-16

Quant Time: Sep 23 09:18:11 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 20 09:45:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	77915	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	129566	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	126198	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	57690	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	78789	55.148	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 110.300%	
35) Dibromofluoromethane	5.391	113	67730	50.583	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.160%	
50) Toluene-d8	8.653	98	265476	51.429	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.860%	
62) 4-Bromofluorobenzene	11.079	95	85718	49.023	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 98.040%	
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	19865	22.736	ug/l	89
20) Methylene Chloride	2.788	84	8674	4.675	ug/l	88
25) 2-Butanone	4.580	43	4463	4.046	ug/l	92
37) Ethyl Acetate	4.721	43	10846	4.414	ug/l	99
40) Benzene	6.043	78	123824	20.153	ug/l	96
43) Isopropyl Acetate	6.342	43	47688	12.629	ug/l	99
52) Toluene	8.720	92	82843	20.674	ug/l	97
67) Ethyl Benzene	10.195	91	9360	1.362	ug/l	96
68) m/p-Xylenes	10.299	106	15175	5.496	ug/l	96
69) o-Xylene	10.640	106	6673	2.532	ug/l	93
70) Styrene	10.659	104	5285	1.218	ug/l	91
84) 1,2,4-Trimethylbenzene	11.756	105	7601	2.111	ug/l	74
95) Naphthalene	13.774	128	76362	18.464	ug/l	100

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

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