

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031522\
 Data File : VX027455.D
 Acq On : 15 Mar 2022 18:19
 Operator : JC/MD
 Sample : N1924-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 33605

Quant Time: Mar 16 04:20:49 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X030822W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 15:54:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	368384	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	596938	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	568519	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	312231	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	222448	51.781	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 103.560%			
35) Dibromofluoromethane	5.391	113	192746	50.040	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 100.080%			
50) Toluene-d8	8.653	98	716354	49.262	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 98.520%			
62) 4-Bromofluorobenzene	11.079	95	297265	53.636	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 107.280%			
Target Compounds						Qvalue
11) Tert butyl alcohol	2.965	59	25720	30.540	ug/l	96
13) Acrolein	2.233	56	5050	8.555	ug/l	85
16) Acetone	2.380	43	161705	95.337	ug/l	95
18) Methyl Acetate	2.709	43	13166	2.788	ug/l	97
25) 2-Butanone	4.568	43	22074	7.460	ug/l	100
40) Benzene	6.044	78	163227	10.021	ug/l	99
52) Toluene	8.720	92	108109	10.411	ug/l	100
64) Tetrachloroethene	9.275	164	71683	17.992	ug/l	99
67) Ethyl Benzene	10.195	91	180294	8.741	ug/l	98
68) m/p-Xylenes	10.299	106	313281	38.721	ug/l	98
69) o-Xylene	10.647	106	84900	10.628	ug/l	99
70) Styrene	10.659	104	72283	5.451	ug/l	89
95) Naphthalene	13.774	128	272985	11.606	ug/l	99

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

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