

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010522\
 Data File : VX026330.D
 Acq On : 05 Jan 2022 18:19
 Operator : JC/MD
 Sample : N1005-08DL 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KEAR-SBDL

Quant Time: Jan 06 03:11:10 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	148543	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	251926	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	219808	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	97244	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	109383	50.900	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 101.800%		
35) Dibromofluoromethane	5.391	113	81418	49.710	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 99.420%		
50) Toluene-d8	8.653	98	294767	48.384	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 96.760%		
62) 4-Bromofluorobenzene	11.079	95	111827	51.563	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 103.120%		
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.971	59	34368	80.070	ug/l	98
16) Acetone	2.386	43	2048	1.759	ug/l	97
31) Cyclohexane	5.476	56	10431	3.408	ug/l	97
39) Methylcyclohexane	7.385	83	11592	3.890	ug/l	93
40) Benzene	6.043	78	902981	119.768	ug/l	97
52) Toluene	8.720	92	59025	12.717	ug/l	98
67) Ethyl Benzene	10.195	91	2709	0.314	ug/l	93
68) m/p-Xylenes	10.305	106	3231	1.002	ug/l	91
69) o-Xylene	10.646	106	1168	0.370	ug/l	92

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

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