

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL110922\
 Data File : VL039857.D
 Acq On : 9 Nov 2022 20:50
 Operator : SY/AP
 Sample : N5527-04DL 10X
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 10605-BASEMENT-NEAR-SINKDL

Quant Time: Nov 10 03:51:52 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL110222AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2022
 QLast Update : Thu Nov 03 09:27:12 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Semsettin Yesilyurt 11/10/2022
 Supervised By : Mahesh Dadoda 11/11/2022

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

Internal Standards						
1) Bromochloromethane	5.179	49	980874	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	6.632	114	2332219	10.000	ppbv	-0.01
55) Chlorobenzene-d5	11.433	117	2086812	10.000	ppbv	-0.02
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	13.744	95	1692361	10.835	ppbv	-0.02
Spiked Amount	10.000	Range	65 - 135	Recovery	= 108.300%	
Target Compounds						
					Qvalue	
3) Chlorodifluoromethane	2.684	51	1047567	8.958	ppbv	97
9) Propene	2.713	41	77843	1.278	ppbv #	65
13) Ethanol	3.240	45	362117	84.096	ppbv	98
15) Acetone	3.468	43	1328091	11.124	ppbv	100
19) Isopropyl Alcohol	3.581	45	140941m	2.385	ppbv	
34) 2-Butanone	4.780	43	186701	1.940	ppbv	93
41) Tetrahydrofuran	5.562	42	46459m	0.578	ppbv	
53) Toluene	9.192	91	140864	0.624	ppbv	97
59) m/p-Xylene	12.317	91	76520m	0.325	ppbv	

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

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