

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL030122\
 Data File : VL038453.D
 Acq On : 1 Mar 2022 16:28
 Operator : SY/AP
 Sample : N1720-01 4X
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 SV1

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 03/02/2022
 Supervised By :Mahesh Dadoda 03/02/2022

Quant Time: Mar 02 03:06:23 2022

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_L\METHODS\VL022222AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Tue Feb 22 2022

QLast Update : Tue Feb 22 19:45:51 2022

Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	5.668	49	902471	10.000	ppbv	0.02
33) 1,4-Difluorobenzene	7.176	114	2282600	10.000	ppbv	0.02
55) Chlorobenzene-d5	12.082	117	2041622	10.000	ppbv	# 0.03
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	14.410	95	1603049	9.490	ppbv	0.02
Spiked Amount	10.000	Range	65 - 135	Recovery	=	94.900%
Target Compounds						Qvalue
2) Dichlorodifluoromethane	3.060	85	16377m	0.115	ppbv	
9) Propene	3.021	41	31165	0.592	ppbv	100
10) Heptane	8.121	43	61347	0.479	ppbv	91
13) Ethanol	3.632	45	20236m	3.854	ppbv	
15) Acetone	3.873	43	146999	1.915	ppbv	# 84
19) Isopropyl Alcohol	4.002	45	13615m	0.306	ppbv	
20) Methylene Chloride	4.349	84	30592	0.805	ppbv	98
26) Hexane	5.693	57	656277	7.029	ppbv	# 27
28) Methyl tert-Butyl Ether	5.060	73	25539m	0.269	ppbv	
29) Chloroform	5.751	83	160949	1.056	ppbv	94
34) 2-Butanone	5.269	43	34855	0.486	ppbv	# 76
52) Tetrachloroethene	11.224	164	876376	13.138	ppbv	98
53) Toluene	9.803	91	42125	0.193	ppbv	91
59) m/p-Xylene	12.963	91	49075m	0.205	ppbv	
60) o-Xylene	13.687	91	24208m	0.107	ppbv	
61) Styrene	13.529	104	23037m	0.220	ppbv	
62) Isopropylbenzene	14.664	105	83709	0.275	ppbv	96

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

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