

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL102422\
 Data File : VL039778.D
 Acq On : 25 Oct 2022 4:49
 Operator : SY/AP
 Sample : N5227-05
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 E

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 10/28/2022
 Supervised By :Mahesh Dadoda 10/31/2022

Quant Time: Oct 25 15:02:38 2022

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL102422AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Mon Oct 24 20:00:55 2022

Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	5.201	49	962073	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	6.658	114	2450472	10.000	ppbv	0.00
55) Chlorobenzene-d5	11.462	117	2123863	10.000	ppbv	0.00
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	13.770	95	1524149	10.088	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	= 100.900%	
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	2.751	85	88529	0.879	ppbv	99
3) Chlorodifluoromethane	2.697	51	729519	6.312	ppbv	96
11) Trichlorofluoromethane	3.581	101	134221	0.954	ppbv	91
12) 1,1,2-Trichlorotrifluo...	4.083	101	27698	0.254	ppbv	94
15) Acetone	3.488	43	471177	4.100	ppbv	96
18) 1,1-Dichloroethene	3.899	96	533625	11.110	ppbv	95
20) Methylene Chloride	3.951	84	183521	4.198	ppbv	96
26) Hexane	5.218	57	135025	1.445	ppbv #	38
29) Chloroform	5.282	83	33665	0.229	ppbv	99
32) 1,1,1-Trichloroethane	6.009	97	464018	3.481	ppbv	97
36) Benzene	6.375	78	68383	0.362	ppbv	96
38) Trichloroethene	7.295	130	2724527	25.574	ppbv	96
52) Tetrachloroethene	10.613	164	151177	2.139	ppbv	97
53) Toluene	9.221	91	74865	0.379	ppbv	91
59) m/p-Xylene	12.336	91	52071m	0.238	ppbv	
60) o-Xylene	13.053	91	23915m	0.117	ppbv	

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

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