

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL102522\
 Data File : VL039802.D
 Acq On : 26 Oct 2022 00:41
 Operator : SY/AP
 Sample : N5227-10
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 J

Manual Integrations
 APPROVED

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

Quant Time: Oct 27 16:31:07 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	5.205	49	838290	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	6.664	114	2261995	10.000	ppbv	0.00
55) Chlorobenzene-d5	11.471	117	1944458	10.000	ppbv	0.00
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	13.777	95	1386830	10.026	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	100.300%
Target Compounds						Qvalue
2) Dichlorodifluoromethane	2.751	85	27216	0.310	ppbv	91
9) Propene	2.716	41	187500	4.334	ppbv	93
11) Trichlorofluoromethane	3.581	101	29379	0.240	ppbv	88
15) Acetone	3.491	43	687743	6.869	ppbv	98
19) Isopropyl Alcohol	3.613	45	31220m	0.665	ppbv	
20) Methylene Chloride	3.954	84	74245	1.949	ppbv	87
26) Hexane	5.227	57	74831	0.919	ppbv #	35
27) Carbon Disulfide	4.144	76	27816	0.288	ppbv #	75
34) 2-Butanone	4.819	43	62790	0.590	ppbv #	88
35) Carbon Tetrachloride	6.500	117	9181m	0.071	ppbv	
36) Benzene	6.378	78	72382m	0.416	ppbv	
53) Toluene	9.224	91	78001	0.427	ppbv	98
58) Ethyl Benzene	12.098	91	21739m	0.098	ppbv	
59) m/p-Xylene	12.355	91	61933m	0.310	ppbv	
60) o-Xylene	13.063	91	28201m	0.150	ppbv	
74) 1,2,4-Trimethylbenzene	16.121	105	62860m	0.287	ppbv	
75) 1,3-Dichlorobenzene	16.326	146	23372m	0.176	ppbv	

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

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