

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL080824\
 Data File : VL041518.D
 Acq On : 8 Aug 2024 18:34
 Operator : SY/MD
 Sample : P3496-05DL 20X
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 SV-4DL

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 08/09/2024
 Supervised By :Mahesh Dadoda 08/09/2024

Quant Time: Aug 09 05:57:51 2024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL072524AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LFri Aug 9

QLast Update : Tue Aug 06 01:32:58 2024

Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	5.185	49	1171233	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	6.648	114	3032376	10.000	ppbv	0.00
55) Chlorobenzene-d5	11.458	117	2752972	10.000	ppbv	# 0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene 13.767 95 2127243 10.041 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 100.400%

Target Compounds

						Qvalue
4) Chloromethane	2.806	50	8531	0.115	ppbv	97
9) Propene	2.690	41	466647	7.286	ppbv	98
10) Heptane	7.571	43	77404	0.538	ppbv	97
13) Ethanol	3.237	45	343058m	59.926	ppbv	
15) Acetone	3.459	43	4542272	34.133	ppbv #	89
17) tert-Butyl alcohol	3.877	59	72396	0.597	ppbv #	94
19) Isopropyl Alcohol	3.575	45	59874	0.770	ppbv #	93
20) Methylene Chloride	3.928	84	91545	1.845	ppbv	95
26) Hexane	5.208	57	69422	0.633	ppbv #	7
27) Carbon Disulfide	4.115	76	39641m	0.309	ppbv	
34) 2-Butanone	4.774	43	6793508	46.313	ppbv	97
51) 2-Hexanone	9.561	43	58925m	0.337	ppbv	
53) Toluene	9.214	91	24629m	0.100	ppbv	
58) Ethyl Benzene	12.076	91	70952	0.224	ppbv	93
59) m/p-Xylene	12.339	91	158942m	0.580	ppbv	
60) o-Xylene	13.047	91	51159	0.197	ppbv	98

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

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