

Data Path : W:\HPCHEM1\MSVOA L\DATA\VL122017\
 Data File : VL031274.D
 Acq On : 20 Dec 2017 17:26
 Operator : SY/AP
 Sample : I7005-01DUP
 Misc : 400ml/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 EP1-OUT-01DUP

Manual Integrations
 APPROVED

MMDadoda
 12/22/2017 4:54:50 PM

Quant Time: Dec 22 04:30:23 2017
 Quant Method : W:\HPCHEM1\MSVOA L\METHODS\VL121317AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Thu Dec 14 2017
 QLast Update : Thu Dec 14 04:29:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.91	49	1494873	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.45	114	3622023	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.41	117	3447469	10.00	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	14.75	95	2718311	11.09	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	110.90%

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	3.18	85	52300	0.41	ppbv	92
3) Chlorodifluoromethane	3.12	51	184381	1.08	ppbv	99
4) Chloromethane	3.27	50	45333	0.49	ppbv	92
9) Propene	3.14	41	74980	0.90	ppbv	85
10) Heptane	8.41	43	28388	0.11	ppbv	100
11) Trichlorofluoromethane	4.13	101	60758	0.29	ppbv	94
13) Ethanol	3.77	45	199098	8.27	ppbv	100
15) Acetone	4.03	43	681764	3.46	ppbv	90
20) Methylene Chloride	4.54	84	577478	8.84	ppbv	96
26) Hexane	5.93	57	1443451	7.55	ppbv #	44
30) Cyclohexane	7.41	84	20651m	0.12	ppbv	
34) 2-Butanone	5.48	43	48440	0.19	ppbv	98
35) Carbon Tetrachloride	7.30	117	18645m	0.07	ppbv	
36) Benzene	7.16	78	111200	0.29	ppbv	98
44) 2,2,4-Trimethylpentane	8.15	57	94021	0.16	ppbv #	82
52) Tetrachloroethene	11.55	164	7285m	0.05	ppbv	
53) Toluene	10.12	91	200600	0.48	ppbv	97
59) m/p-Xylene	13.30	91	97813m	0.21	ppbv	

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

