

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL111620\
 Data File : VL035962.D
 Acq On : 17 Nov 2020 3:07
 Operator : SY/AP
 Sample : L4773-01 10X
 Misc : 400ml/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 SV1

Manual Integrations
 APPROVED

MMDadoda
 11/18/2020 9:08:32 AM

Quant Time: Nov 17 22:29:08 2020
 Quant Title :
 QLast Update : Wed Nov 04 02:40:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.67	49	1135838	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.18	114	4025618	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.09	117	4101806	10.00	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	14.42	95	3156018	10.83	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	108.30%

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	3.04	85	22038	0.103	ppbv	95
3) Chlorodifluoromethane	2.98	51	22986	0.169	ppbv	93
9) Propene	3.01	41	11109	0.223	ppbv	96
11) Trichlorofluoromethane	3.95	101	24333	0.091	ppbv	100
13) Ethanol	3.60	45	482217	44.647	ppbv	97
15) Acetone	3.85	43	734543	6.049	ppbv	97
19) Isopropyl Alcohol	3.98	45	39068m	0.524	ppbv	
29) Chloroform	5.75	83	1163853	4.488	ppbv	97
31) cis-1,2-Dichloroethene	5.55	61	15866m	0.117	ppbv	
38) Trichloroethene	7.83	130	1118835	7.654	ppbv	97
52) Tetrachloroethene	11.23	164	11958387m	84.340	ppbv	
59) m/p-Xylene	12.98	91	64168m	0.152	ppbv	
60) o-Xylene	13.70	91	44300	0.108	ppbv	87

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

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