

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF010919\
 Data File : VF061302.D
 Acq On : 9 Jan 2019 15:42
 Operator : VA/AP
 Sample : K1082-09
 Misc : 4.67µ/5mL/MSVOA-F/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 SP-1-5

Quant Time: Jan 10 05:55:13 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F010819S.M
 Quant Title : SW846 8260
 QLast Update : Wed Jan 09 07:13:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	0.00	168	0	0.00	ug/l	-4.87
34) 1,4-Difluorobenzene	0.00	114	0	0.00	ug/l	-5.60
63) Chlorobenzene-d5	0.00	117	0	0.00	ug/l	-9.79
72) 1,4-Dichlorobenzene-d4	12.50	152	67	50.00	ug/l	-0.05

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	0.00	65	0	0.00	ug/l	
Spiked Amount	50.000		Recovery	=	0.00%	
35) Dibromofluoromethane	0.00	113	0	0.00	ug/l	
Spiked Amount	50.000		Recovery	=	0.00%	
50) Toluene-d8	0.00	98	0	0.00	ug/l	
Spiked Amount	50.000		Recovery	=	0.00%	
62) 4-Bromofluorobenzene	0.00	95	0	0.00	ug/l	
Spiked Amount	50.000		Recovery	=	0.00%	

Target Compounds

						Qvalue
74) N-amyl acetate	11.36	43	459	342.490	ug/l #	48
75) 1,1,2,2-Tetrachloroethane	11.65	83	69	73.813	ug/l #	25
78) n-propylbenzene	11.71	91	94	13.372	ug/l #	55
79) 2-Chlorotoluene	11.71	91	94	24.002	ug/l #	41
81) trans-1,4-Dichloro-2-buten	11.96	75	91	303.730	ug/l #	17
89) n-Butylbenzene	12.71	91	77	13.845	ug/l #	27
92) 1,2-Dibromo-3-Chloropropan	13.52	75	131	829.405	ug/l #	1

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

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