

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF031119\
 Data File : VF061896.D
 Acq On : 11 Mar 2019 18:24
 Operator : VA/AP
 Sample : K1824-17
 Misc : 5.04g/5mL/MSVOA-F/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 P-3

Quant Time: Mar 12 05:13:17 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F030419S.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 05 04:50:44 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	-4.88
34) 1,4-Difluorobenzene	0.00	114	0	0.00	ug/l	-5.61
63) Chlorobenzene-d5	0.00	117	0	0.00	ug/l	-9.78
72) 1,4-Dichlorobenzene-d4	12.55	152	71	50.00	ug/l	0.00
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	7.53	98	64	0.00	ug/l	-0.03
Spiked Amount	50.000		Recovery	=	0.00%	
62) 4-Bromofluorobenzene	11.48	95	63	0.00	ug/l	0.06
Spiked Amount	50.000		Recovery	=	0.00%	
Target Compounds						
					Qvalue	
74) N-amyl acetate	11.39	43	70	62.207	ug/l #	42
80) 1,3,5-Trimethylbenzene	11.78	105	63	16.053	ug/l #	24
83) tert-Butylbenzene	12.13	119	66	15.838	ug/l #	36
89) n-Butylbenzene	12.81	91	97	22.019	ug/l #	26
90) Hexachloroethane	12.88	117	110	99.143	ug/l #	24
92) 1,2-Dibromo-3-Chloropropan	13.70	75	85	560.896	ug/l #	1

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

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