

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX043019\
 Data File : VX009241.D
 Acq On : 30 Apr 2019 13:20
 Operator : JC/SP
 Sample : K2618-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CARBON-DNM

Quant Time: May 01 02:08:32 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 30 07:03:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	182079	50.00	ug/l	-0.02
34) 1,4-Difluorobenzene	6.85	114	290424	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	258431	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	113705	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	110599	44.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.00%	
35) Dibromofluoromethane	5.52	113	68441	37.49	ug/l	0.02
Spiked Amount 50.000			Recovery	=	74.98%	
50) Toluene-d8	8.71	98	371156	50.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.70%	
62) 4-Bromofluorobenzene	11.14	95	124212	46.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.76%	
Target Compounds						
3) Chloromethane	1.33	50	40479	17.671	ug/l	97
16) Acetone	2.53	43	99779	74.759	ug/l	98
18) Methyl Acetate	2.82	43	699692	220.437	ug/l	98
29) Tetrahydrofuran	5.22	42	108019	80.935	ug/l	94

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

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