

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041319\
 Data File : VX008929.D
 Acq On : 13 Apr 2019 14:12
 Operator : JC/SP
 Sample : K2371-01DL 100X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OK-COMPDL

Quant Time: Apr 15 02:26:33 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X040319W.M
 Quant Title : SW846 8260
 QLast Update : Thu Apr 04 04:47:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	196692	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	315309	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	268320	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	115672	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	126177	43.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.96%	
35) Dibromofluoromethane	5.50	113	93785	46.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.04%	
50) Toluene-d8	8.71	98	386659	48.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.34%	
62) 4-Bromofluorobenzene	11.13	95	129313	45.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.86%	
Target Compounds						
16) Acetone	2.44	43	97801	59.188	ug/l	99
52) Toluene	8.79	92	137062	22.642	ug/l	99
67) Ethyl Benzene	10.25	91	6774	0.611	ug/l	98
68) m/p-Xylenes	10.36	106	10381	2.603	ug/l	98
69) o-Xylene	10.69	106	4523	1.170	ug/l	92
84) 1,2,4-Trimethylbenzene	11.80	105	6923	0.850	ug/l	98

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

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