

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041919\
 Data File : VX009064.D
 Acq On : 19 Apr 2019 12:46
 Operator : JC/SP
 Sample : K2482-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 AU-041819

Quant Time: Apr 22 11:08:45 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X041519W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 17 06:47:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	229684	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	373031	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	319655	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	138901	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.06	65	155553	48.06	ug/l	0.00
Spiked Amount			Recovery	=	96.12%	
35) Dibromofluoromethane	5.50	113	113940	47.59	ug/l	0.00
Spiked Amount			Recovery	=	95.18%	
50) Toluene-d8	8.71	98	465785	48.79	ug/l	0.00
Spiked Amount			Recovery	=	97.58%	
62) 4-Bromofluorobenzene	11.13	95	155406	46.11	ug/l	0.00
Spiked Amount			Recovery	=	92.22%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.45	43	920499	548.379	ug/l	99
17) Carbon Disulfide	2.57	76	2746	1.623	ug/l	99
25) 2-Butanone	4.68	43	1097253	408.950	ug/l	99
30) Chloroform	5.22	83	2790	0.595	ug/l	90
51) 4-Methyl-2-Pentanone	8.64	43	75965	14.905	ug/l	100
52) Toluene	8.79	92	2191202	321.245	ug/l	98
68) m/p-Xylenes	10.36	106	4171	0.931	ug/l	95
69) o-Xylene	10.70	106	3572	0.819	ug/l	98
84) 1,2,4-Trimethylbenzene	11.80	105	3034	0.337	ug/l	87
95) Naphthalene	13.83	128	4660	0.439	ug/l	99

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

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