

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX051920\
 Data File : VX016343.D
 Acq On : 19 May 2020 18:27
 Operator : JC/SP
 Sample : L2663-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1366

Quant Time: May 20 07:33:39 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X050420W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 06 06:43:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	225939	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	378100	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	366794	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	180507	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.03	65	134226	45.15	ug/l	0.00
Spiked Amount			Recovery	=	90.30%	
35) Dibromofluoromethane	5.47	113	110304	46.05	ug/l	0.00
Spiked Amount			Recovery	=	92.10%	
50) Toluene-d8	8.70	98	460351	49.88	ug/l	0.00
Spiked Amount			Recovery	=	99.76%	
62) 4-Bromofluorobenzene	11.12	95	184067	52.43	ug/l	0.00
Spiked Amount			Recovery	=	104.86%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
11) Tert butyl alcohol	3.01	59	24519	32.613	ug/l	98
16) Acetone	2.42	43	1008789	760.443	ug/l	94
17) Carbon Disulfide	2.55	76	5744	0.610	ug/l #	88
18) Methyl Acetate	2.75	43	2142	0.657	ug/l	94
25) 2-Butanone	4.64	43	232435	108.733	ug/l	97
29) Tetrahydrofuran	5.10	42	2339	1.651	ug/l #	91
30) Chloroform	5.18	83	1348	0.298	ug/l	99
40) Benzene	6.12	78	9814	0.901	ug/l #	89
51) 4-Methyl-2-Pentanone	8.62	43	34538	8.200	ug/l	96
52) Toluene	8.77	92	5190	0.765	ug/l	97
59) 2-Hexanone	9.48	43	47862	14.470	ug/l	93
68) m/p-Xylenes	10.34	106	2349	0.458	ug/l	92
84) 1,2,4-Trimethylbenzene	11.79	105	5473	0.488	ug/l #	60
86) p-Isopropyltoluene	12.05	119	18996	1.586	ug/l	95

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

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