

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX050518\
 Data File : VX001433.D
 Acq On : 05 May 2018 21:55
 Operator : JC/MD
 Sample : J2695-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CS-DRUM-2

Quant Time: May 06 08:27:22 2018
 Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X050518W.M
 Quant Title : SW846 8260
 QLast Update : Sun May 06 07:59:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	243341	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	329698	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	290644	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.09	152	165729	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.07	65	160119	45.78	ug/l	0.00
Spiked Amount			Recovery	=	91.56%	
35) Dibromofluoromethane	5.51	113	123486	42.11	ug/l	0.00
Spiked Amount			Recovery	=	84.22%	
50) Toluene-d8	8.72	98	462268	43.18	ug/l	0.00
Spiked Amount			Recovery	=	86.36%	
62) 4-Bromofluorobenzene	11.14	95	155727	37.90	ug/l	0.00
Spiked Amount			Recovery	=	75.80%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
11) Tert butyl alcohol	3.08	59	6708	9.097	ug/l #	94
16) Acetone	2.45	43	682308	429.617	ug/l	99
17) Carbon Disulfide	2.57	76	14256	2.080	ug/l	99
25) 2-Butanone	4.72	43	266898	116.258	ug/l	97
30) Chloroform	5.23	83	16105	2.984	ug/l	96
40) Benzene	6.15	78	4734	0.386	ug/l	98
51) 4-Methyl-2-Pentanone	8.66	43	245050	52.253	ug/l	97
52) Toluene	8.79	92	9923	1.239	ug/l	99
59) 2-Hexanone	9.51	43	41538	11.567	ug/l	100
67) Ethyl Benzene	10.26	91	3825	0.274	ug/l	98
68) m/p-Xylenes	10.36	106	5864	1.062	ug/l	98
69) o-Xylene	10.70	106	2790	0.520	ug/l	97
84) 1,2,4-Trimethylbenzene	11.81	105	10687	1.089	ug/l	79
86) p-Isopropyltoluene	12.07	119	3324	0.318	ug/l #	71
95) Naphthalene	13.84	128	5878	0.479	ug/l	95

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

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