

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061120\
 Data File : VX016713.D
 Acq On : 11 Jun 2020 18:36
 Operator : JC/SP
 Sample : L2977-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 I-OC-2

Quant Time: Jun 12 07:49:54 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X052720W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 27 16:31:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	206728	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	327282	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	331518	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	180133	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	122106	47.56	ug/l	0.00
Spiked Amount			Recovery	=	95.12%	
35) Dibromofluoromethane	5.47	113	102188	50.71	ug/l	0.00
Spiked Amount			Recovery	=	101.42%	
50) Toluene-d8	8.70	98	409111	52.16	ug/l	0.00
Spiked Amount			Recovery	=	104.32%	
62) 4-Bromofluorobenzene	11.12	95	166624	54.98	ug/l	0.00
Spiked Amount			Recovery	=	109.96%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.42	43	457370	406.627	ug/l	99
17) Carbon Disulfide	2.55	76	540997	91.584	ug/l	99
25) 2-Butanone	4.64	43	83915	44.029	ug/l	94
30) Chloroform	5.18	83	151494	37.906	ug/l	100
40) Benzene	6.12	78	134914	14.620	ug/l	98
47) Bromodichloromethane	7.88	83	9619	2.944	ug/l #	97
52) Toluene	8.77	92	36929	6.346	ug/l	96
67) Ethyl Benzene	10.23	91	36586	3.040	ug/l	100
68) m/p-Xylenes	10.34	106	52635	11.797	ug/l	100
69) o-Xylene	10.68	106	24213	5.604	ug/l	98
84) 1,2,4-Trimethylbenzene	11.79	105	34251	3.250	ug/l	93
95) Naphthalene	13.82	128	3134129	250.531	ug/l	99

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

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