

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051421\
 Data File : VX021987.D
 Acq On : 14 May 2021 12:01
 Operator : JC/MD
 Sample : M2383-08DL 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KTW-02DL

Manual Integrations
 APPROVED

MMDadoda
 5/17/2021 2:34:17 PM

Quant Time: May 15 01:03:01 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050721W.M
 Quant Title : SW846 8260
 QLast Update : Sat May 08 05:38:18 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.562	168	78662	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	170653	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	182030	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	85533	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	66118	51.07	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 102.14%			
35) Dibromofluoromethane	5.391	113	57713	49.35	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 98.70%			
50) Toluene-d8	8.653	98	225400	50.50	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 101.00%			
62) 4-Bromofluorobenzene	11.085	95	101319	57.52	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 115.04%			
Target Compounds					Qvalue	
16) Acetone	2.386	43	1277	2.20	ug/l	90
17) Carbon Disulfide	2.514	76	272087	91.72	ug/l	99
30) Chloroform	5.105	83	1464m	0.69	ug/l	
40) Benzene	6.050	78	19094	3.51	ug/l #	88
52) Toluene	8.726	92	3814	1.10	ug/l	98
67) Ethyl Benzene	10.195	91	27382	3.52	ug/l	98
68) m/p-Xylenes	10.305	106	18117	6.10	ug/l	99
69) o-Xylene	10.646	106	6907	2.33	ug/l	94
73) Isopropylbenzene	10.963	105	1957	0.25	ug/l	93
78) n-propylbenzene	11.311	91	8004	0.87	ug/l	96
80) 1,3,5-Trimethylbenzene	11.457	105	15373	2.31	ug/l	98
84) 1,2,4-Trimethylbenzene	11.756	105	55689	8.35	ug/l	100
95) Naphthalene	13.780	128	444088	57.92	ug/l	99

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

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