

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN071318\
 Data File : VN049793.D
 Acq On : 13 Jul 2018 13:32
 Operator : MD\SY
 Sample : J3762-09 MDL 0.5ppb
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MDL-WATER-03-QT3-2018

Manual Integrations
 APPROVED

MMDadoda
 7/16/2018 11:25:26 AM

Quant Time: Jul 16 04:00:58 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N071218W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 12 15:51:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1069895	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1688634	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1450626	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	555599	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	690931	55.51	ug/l	0.00
Spiked Amount	50.000		Recovery	=	111.02%	
35) Dibromofluoromethane	7.59	113	657061	51.62	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.24%	
50) Toluene-d8	10.09	98	1973867	45.30	ug/l	0.00
Spiked Amount	50.000		Recovery	=	90.60%	
62) 4-Bromofluorobenzene	12.40	95	599412	38.52	ug/l	0.00
Spiked Amount	50.000		Recovery	=	77.04%	
Target Compounds						
						Qvalue
2) Dichlorodifluoromethane	1.85	85	7111	Below Cal		84
3) Chloromethane	2.06	50	10786	Below Cal		98
5) Bromomethane	2.57	94	9746	Below Cal		83
16) Acetone	3.83	43	11757	1.78	ug/l	90
17) Carbon Disulfide	4.05	76	23980	0.38	ug/l	100
41) Methacrylonitrile	7.18	41	2758	0.42	ug/l #	75
49) 1,4-Dioxane	9.20	88	1603	11.82	ug/l #	53
75) 1,1,2,2-Tetrachloroethane	12.51	83	9123	Below Cal		92
81) trans-1,4-Dichloro-2-buten	12.30	75	1705m	0.56	ug/l	
92) 1,2-Dibromo-3-Chloropropan	14.27	75	1356	0.35	ug/l	75
94) Hexachlorobutadiene	15.01	225	4695	Below Cal		96
96) 1,2,3-Trichlorobenzene	15.31	180	4976	0.50	ug/l	93

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

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