

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN060520\
 Data File : VN061666.D
 Acq On : 5 Jun 2020 13:57
 Operator : JC/MD
 Sample : L2904-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VE-3-1

Manual Integrations
 APPROVED

MMDadoda
 6/8/2020 11:49:07 AM

Quant Time: Jun 06 04:36:20 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N052820W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 29 02:14:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	294753	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	513107	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	506530	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	246808	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	171855	48.94	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.88%	
35) Dibromofluoromethane	7.55	113	146694	56.36	ug/l	0.00
Spiked Amount	50.000		Recovery	=	112.72%	
50) Toluene-d8	10.06	98	615733	51.05	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.10%	
62) 4-Bromofluorobenzene	12.38	95	227352	51.66	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.32%	
Target Compounds						
						Qvalue
16) Acetone	3.78	43	31511	27.027	ug/l	100
25) 2-Butanone	6.79	43	13169	7.912	ug/l	96
65) Chlorobenzene	11.41	112	24749	2.351	ug/l	96
67) Ethyl Benzene	11.48	91	5965	0.328	ug/l	98
69) o-Xylene	11.92	106	6213	0.907	ug/l	95
73) Isopropylbenzene	12.22	105	4975	0.276	ug/l	96
78) n-propylbenzene	12.57	91	9855	0.490	ug/l	99
80) 1,3,5-Trimethylbenzene	12.71	105	28450	1.890	ug/l	97
84) 1,2,4-Trimethylbenzene	13.01	105	54448	3.640	ug/l	99
85) sec-Butylbenzene	13.15	105	7930	0.463	ug/l	92
86) p-Isopropyltoluene	13.26	119	10491	0.665	ug/l	95
87) 1,3-Dichlorobenzene	13.26	146	4234	0.520	ug/l	91
88) 1,4-Dichlorobenzene	13.34	146	13861m	1.671	ug/l	
91) 1,2-Dichlorobenzene	13.63	146	17884	2.298	ug/l	89

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

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