

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN050420\
 Data File : VN061281.D
 Acq On : 4 May 2020 16:26
 Operator : JC/MD
 Sample : L2457-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1341

Manual Integrations
 APPROVED

MMDadoda
 5/5/2020 11:36:20 AM

Quant Time: May 05 05:02:37 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042720W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 29 08:14:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	105202	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	191378	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	181227	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	84790	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	86311	51.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.70%	
35) Dibromofluoromethane	7.55	113	55715	49.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.46%	
50) Toluene-d8	10.06	98	236126	51.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.36%	
62) 4-Bromofluorobenzene	12.38	95	91514	52.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.42%	
Target Compounds						
11) Tert butyl alcohol	4.73	59	16488	73.941	ug/l #	86
16) Acetone	3.77	43	54355	80.688	ug/l	98
25) 2-Butanone	6.78	43	34357	34.880	ug/l	96
40) Benzene	8.00	78	1914	0.335	ug/l #	82
51) 4-Methyl-2-Pentanone	9.95	43	11911	5.429	ug/l	91
70) Styrene	11.94	104	2966	0.749	ug/l	89
86) p-Isopropyltoluene	13.26	119	10674m	1.998	ug/l	

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

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