

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102720\
 Data File : VN064333.D
 Acq On : 27 Oct 2020 18:02
 Operator : JC/MD
 Sample : L4486-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1836

Quant Time: Oct 28 04:48:40 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N102720W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 27 12:23:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	220012	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	382059	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	379110	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	168678	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	7.99	65	181095	51.22	ug/l	0.00
Spiked Amount			Recovery	=	102.44%	
35) Dibromofluoromethane	7.55	113	130318	51.15	ug/l	0.00
Spiked Amount			Recovery	=	102.30%	
50) Toluene-d8	10.06	98	480708	51.01	ug/l	0.00
Spiked Amount			Recovery	=	102.02%	
62) 4-Bromofluorobenzene	12.38	95	206248	56.13	ug/l	0.00
Spiked Amount			Recovery	=	112.26%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
11) Tert butyl alcohol	4.74	59	13758	26.806	ug/l #	77
16) Acetone	3.78	43	46831	32.460	ug/l	96
17) Carbon Disulfide	4.01	76	6808	1.095	ug/l	97
25) 2-Butanone	6.80	43	12918	7.277	ug/l #	81
51) 4-Methyl-2-Pentanone	9.95	43	15619	4.036	ug/l	94
52) Toluene	10.13	92	2920	0.431	ug/l	90
68) m/p-Xylenes	11.59	106	2219	0.419	ug/l	89
69) o-Xylene	11.92	106	1784	0.350	ug/l	98
84) 1,2,4-Trimethylbenzene	13.02	105	6437	0.603	ug/l	99
85) sec-Butylbenzene	13.14	105	3737	0.329	ug/l	98
95) Naphthalene	15.11	128	6705	3.062	ug/l #	88

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

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