

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031820\
 Data File : VN060602.D
 Acq On : 18 Mar 2020 19:05
 Operator : JC/MD
 Sample : L1946-04DL 5X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMPDL

Quant Time: Mar 19 07:38:19 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 18 08:49:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	205008	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	341863	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	316796	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	145207	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	8.00	65	130887	46.51	ug/l	0.00
Spiked Amount			Recovery	=	93.02%	
35) Dibromofluoromethane	7.56	113	101287	49.00	ug/l	0.00
Spiked Amount			Recovery	=	98.00%	
50) Toluene-d8	10.08	98	416995	49.39	ug/l	0.00
Spiked Amount			Recovery	=	98.78%	
62) 4-Bromofluorobenzene	12.40	95	156050	50.43	ug/l	0.00
Spiked Amount			Recovery	=	100.86%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.78	43	87766	88.041	ug/l	100
18) Methyl Acetate	4.29	43	23749	8.464	ug/l	99
20) Methylene Chloride	4.52	84	23071	9.408	ug/l	97
25) 2-Butanone	6.81	43	21188	13.135	ug/l	98
51) 4-Methyl-2-Pentanone	9.97	43	10522	3.135	ug/l	97
52) Toluene	10.14	92	181366	30.052	ug/l	99
67) Ethyl Benzene	11.50	91	15129	1.295	ug/l	99
68) m/p-Xylenes	11.61	106	28733	6.621	ug/l	98
69) o-Xylene	11.94	106	12587	3.038	ug/l	92

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

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