

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN040220\
 Data File : VN060833.D
 Acq On : 2 Apr 2020 19:43
 Operator : JC/MD
 Sample : L2104-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHASSIE-WASH

Quant Time: Apr 03 09:25:53 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	205701	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	340319	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	314401	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	145580	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	129192	45.75	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.50%	
35) Dibromofluoromethane	7.56	113	100525	48.85	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.70%	
50) Toluene-d8	10.08	98	417842	49.71	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.42%	
62) 4-Bromofluorobenzene	12.40	95	154261	50.08	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.16%	
Target Compounds						
16) Acetone	3.78	43	34278	34.269	ug/l	98
20) Methylene Chloride	4.53	84	4384	1.782	ug/l	91
25) 2-Butanone	6.81	43	14511	8.965	ug/l	95
43) Isopropyl Acetate	8.14	43	32513	5.537	ug/l #	91
51) 4-Methyl-2-Pentanone	9.97	43	69771	20.880	ug/l	97
52) Toluene	10.14	92	312045	51.940	ug/l	99

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

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