

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080320\
 Data File : VN062734.D
 Acq On : 3 Aug 2020 14:37
 Operator : JC/MD
 Sample : L3539-01 10X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 I-OC-5

Manual Integrations
 APPROVED

MMDadoda
 8/4/2020 1:52:15 PM

Quant Time: Aug 04 07:48:54 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 21 18:48:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	127533	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	224678	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	225877	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	89041	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	110168	55.61	ug/l	0.00
Spiked Amount			Recovery	=	111.22%	
35) Dibromofluoromethane	7.55	113	74632	50.57	ug/l	0.00
Spiked Amount			Recovery	=	101.14%	
50) Toluene-d8	10.06	98	296177	50.30	ug/l	0.00
Spiked Amount			Recovery	=	100.60%	
62) 4-Bromofluorobenzene	12.38	95	110781	51.45	ug/l	0.00
Spiked Amount			Recovery	=	102.90%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.78	43	27899	37.422	ug/l	98
17) Carbon Disulfide	4.01	76	358095	79.812	ug/l	99
25) 2-Butanone	6.80	43	4938m	4.349	ug/l	
30) Chloroform	7.33	83	8947	2.976	ug/l	92
40) Benzene	8.00	78	30375	4.190	ug/l	94
52) Toluene	10.13	92	5519	1.311	ug/l	95
67) Ethyl Benzene	11.48	91	2285	0.263	ug/l #	83
68) m/p-Xylenes	11.59	106	4435	1.421	ug/l	90
69) o-Xylene	11.92	106	1704	0.575	ug/l	97
84) 1,2,4-Trimethylbenzene	13.02	105	2577	0.450	ug/l	97
95) Naphthalene	15.10	128	391793	95.262	ug/l	99

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

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