

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062263.D
 Acq On : 1 Jul 2020 1:54
 Operator : JC/MD
 Sample : L3148-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EPEC-POLYMERS-02

Quant Time: Jul 01 05:51:35 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 24 15:33:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	194068	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	390038	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	395696	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	142915	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	155315	57.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.06%	
35) Dibromofluoromethane	7.55	113	125758	50.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.52%	
50) Toluene-d8	10.06	98	524707	52.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.92%	
62) 4-Bromofluorobenzene	12.38	95	170659	50.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.00%	
Target Compounds						
16) Acetone	3.78	43	19326	18.862	ug/l	91
18) Methyl Acetate	4.28	43	3276	1.513	ug/l #	80
20) Methylene Chloride	4.51	84	22877	8.973	ug/l	94
25) 2-Butanone	6.79	43	8280	6.220	ug/l	98
43) Isopropyl Acetate	8.13	43	49536	9.360	ug/l	98
79) 2-Chlorotoluene	12.65	91	90895	11.882	ug/l	97

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

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