

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062262.D
 Acq On : 1 Jul 2020 1:29
 Operator : JC/MD
 Sample : L3148-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EPEC-POLYMERS-01

Quant Time: Jul 01 05:49:09 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	208622	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	408732	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	411958	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	144937	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	162656	55.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.12%	
35) Dibromofluoromethane	7.55	113	131202	50.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.06%	
50) Toluene-d8	10.06	98	544670	51.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.92%	
62) 4-Bromofluorobenzene	12.38	95	177042	49.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.98%	
Target Compounds						
16) Acetone	3.78	43	32278	31.586	ug/l	98
20) Methylene Chloride	4.51	84	21922	7.975	ug/l	95
25) 2-Butanone	6.80	43	12232	8.548	ug/l	99
43) Isopropyl Acetate	8.13	43	45664	8.234	ug/l	97

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

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