

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062264.D
 Acq On : 1 Jul 2020 2:18
 Operator : JC/MD
 Sample : L3148-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EPEC-POLYMERS-03

Quant Time: Jul 01 05:53:27 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	189196	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	386094	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	382974	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	135258	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	152293	57.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.72%	
35) Dibromofluoromethane	7.55	113	123200	49.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.48%	
50) Toluene-d8	10.06	98	505612	51.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.14%	
62) 4-Bromofluorobenzene	12.38	95	159424	47.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.32%	
Target Compounds						
16) Acetone	3.78	43	33423	36.648	ug/l	98
17) Carbon Disulfide	4.01	76	9474	1.572	ug/l	98
18) Methyl Acetate	4.28	43	3453	1.636	ug/l #	89
20) Methylene Chloride	4.50	84	63164	25.810	ug/l	97
43) Isopropyl Acetate	8.13	43	43644	8.331	ug/l	100
79) 2-Chlorotoluene	12.65	91	41437	5.723	ug/l	97
82) 4-Chlorotoluene	12.75	91	33361	4.456	ug/l	97

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

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