

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN041122\
 Data File : VN071944.D
 Acq On : 11 Apr 2022 18:29
 Operator : JC\MD
 Sample : PB143992ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB143992ZHE#06

Quant Time: Apr 12 08:42:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.081	168	694778	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1204656	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1271126	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	507103	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	383598	48.363	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	96.720%
35) Dibromofluoromethane	8.016	113	340848	48.083	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	96.160%
50) Toluene-d8	10.439	98	1525671	52.612	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	105.220%
62) 4-Bromofluorobenzene	12.727	95	544849	58.096	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	116.200%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	16949	5.000	ug/l #	88
18) Methyl Acetate	4.863	43	22253	1.173	ug/l	97
43) Isopropyl Acetate	8.563	43	190958	10.596	ug/l #	70

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN041122\
Data File : VN071944.D
Acq On : 11 Apr 2022 18:29
Operator : JC\MD
Sample : PB143992ZHE#06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB143992ZHE#06

Quant Time: Apr 12 08:42:03 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260
QLast Update : Thu Mar 31 01:57:54 2022
Response via : Initial Calibration

