

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
 Data File : VN073561.D
 Acq On : 21 Jul 2022 17:34
 Operator : JC\MD
 Sample : PB146347TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB146347TB

Quant Time: Jul 22 08:01:35 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 01:54:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	299020	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	484781	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	464149	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	218702	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	176489	40.885	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	81.780%
35) Dibromofluoromethane	7.951	113	155871	48.327	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.660%
50) Toluene-d8	10.381	98	519864	45.382	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.760%
62) 4-Bromofluorobenzene	12.675	95	219505	49.863	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.720%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	64922	26.794	ug/l	90
20) Methylene Chloride	5.016	84	27868	6.544	ug/l #	87
37) Ethyl Acetate	7.340	43	18368	2.449	ug/l #	95
43) Isopropyl Acetate	8.492	43	53381	4.953	ug/l #	87

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072122\
Data File : VN073561.D
Acq On : 21 Jul 2022 17:34
Operator : JC\MD
Sample : PB146347TB
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB146347TB

Quant Time: Jul 22 08:01:35 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 29 01:54:03 2022
Response via : Initial Calibration

