

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX120120\
 Data File : VX019693.D
 Acq On : 01 Dec 2020 14:24
 Operator : JC/MD
 Sample : PB133320ZHE#06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB133320ZHE#06

Quant Time: Dec 02 03:00:12 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 21 01:57:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	290378	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	483093	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	423831	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	179062	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	6.03	65	185391	49.24	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	98.48%		
35) Dibromofluoromethane	5.46	113	146262	48.23	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.46%		
50) Toluene-d8	8.70	98	580434	49.26	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	98.52%		
62) 4-Bromofluorobenzene	11.12	95	193911	45.74	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	91.48%		

Target Compounds

					Qvalue
16) Acetone	2.43	43	11094	7.620	ug/l 98
18) Methyl Acetate	2.75	43	2445	1.311	ug/l 99
20) Methylene Chloride	2.83	84	22728	6.245	ug/l 98
43) Isopropyl Acetate	6.42	43	32964	4.612	ug/l 99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120120\
Data File : VX019693.D
Acq On : 01 Dec 2020 14:24
Operator : JC/MD
Sample : PB133320ZHE#06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB133320ZHE#06

Quant Time: Dec 02 03:00:12 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260
QLast Update : Sat Nov 21 01:57:45 2020
Response via : Initial Calibration

