

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110722\
 Data File : VN075197.D
 Acq On : 07 Nov 2022 18:59
 Operator : JC\MD
 Sample : N5491-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3

Quant Time: Nov 08 01:13:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 09:48:49 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/08/2022
 Supervised By : Mahesh Dadoda 11/08/2022

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	93685	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	197868	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	176744	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	68098	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	77064	61.051	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 122.100%			
35) Dibromofluoromethane	8.171	113	62382	53.704	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 107.400%			
50) Toluene-d8	10.571	98	231682	51.409	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 102.820%			
62) 4-Bromofluorobenzene	12.853	95	75770	47.862	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 95.720%			
Target Compounds						
16) Acetone	4.441	43	9554	17.875	ug/l #	83
18) Methyl Acetate	5.041	43	4818	2.760	ug/l #	82
20) Methylene Chloride	5.283	84	4540	3.810	ug/l	98
37) Ethyl Acetate	7.571	43	17645	8.947	ug/l #	92
43) Isopropyl Acetate	8.682	43	143307m	46.845	ug/l	

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

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