

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081822\
 Data File : VN073997.D
 Acq On : 18 Aug 2022 12:07
 Operator : JC\MD
 Sample : N4241-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TRE-1630

Quant Time: Aug 18 21:48:12 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081522W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 16 00:45:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	739845	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	1094917	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	1068081	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	587470	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	341897	47.446	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	94.900%
35) Dibromofluoromethane	7.951	113	331886	51.255	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.500%
50) Toluene-d8	10.381	98	1180357	47.067	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.140%
62) 4-Bromofluorobenzene	12.669	95	506885	60.902	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	121.800%
Target Compounds						
					Qvalue	
3) Chloromethane	2.263	50	1737	0.266	ug/l	91
5) Bromomethane	2.834	94	2271	0.434	ug/l #	70
11) Tert butyl alcohol	5.281	59	75821	47.039	ug/l	99
13) Acrolein	3.981	56	14982	13.279	ug/l	99
16) Acetone	4.216	43	11459299	3370.651	ug/l	95
17) Carbon Disulfide	4.446	76	8052	0.580	ug/l	99
18) Methyl Acetate	4.781	43	57807	4.488	ug/l	98
20) Methylene Chloride	5.016	84	39831	5.552	ug/l	96
24) 1,1-Dichloroethane	6.304	63	11870	1.040	ug/l	95
25) 2-Butanone	7.257	43	253208	46.972	ug/l	99
30) Chloroform	7.745	83	6783	0.531	ug/l #	72
37) Ethyl Acetate	7.334	43	922665	84.679	ug/l	96
40) Benzene	8.387	78	65196	2.297	ug/l	99
42) 1,2-Dichloroethane	8.463	62	4964	0.509	ug/l	87
44) Trichloroethene	9.151	130	8055	0.982	ug/l	93
51) 4-Methyl-2-Pentanone	10.269	43	190007	18.105	ug/l	99
52) Toluene	10.445	92	196439	10.092	ug/l	98
67) Ethyl Benzene	11.786	91	88159	2.335	ug/l	100
68) m/p-Xylenes	11.886	106	148796	9.701	ug/l	100
69) o-Xylene	12.216	106	112594	7.320	ug/l	98
73) Isopropylbenzene	12.522	105	22408	0.524	ug/l	97
78) n-propylbenzene	12.863	91	50935	1.117	ug/l	96
80) 1,3,5-Trimethylbenzene	12.998	105	97313	2.797	ug/l	100
83) tert-Butylbenzene	13.263	119	10608	0.331	ug/l	95
84) 1,2,4-Trimethylbenzene	13.310	105	577344	16.616	ug/l	80
85) sec-Butylbenzene	13.439	105	26044	0.628	ug/l	89
86) p-Isopropyltoluene	13.557	119	35352	1.000	ug/l	94
89) n-Butylbenzene	13.880	91	29550	1.055	ug/l #	74
95) Naphthalene	15.433	128	600620	14.598	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081822\
Data File : VN073997.D
Acq On : 18 Aug 2022 12:07
Operator : JC\MD
Sample : N4241-07
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TRE-1630

Quant Time: Aug 18 21:48:12 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081522W.M
Quant Title : SW846 8260
QLast Update : Tue Aug 16 00:45:21 2022
Response via : Initial Calibration

