

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092222\
 Data File : VX031585.D
 Acq On : 22 Sep 2022 21:01
 Operator : JC/MD
 Sample : N4642-15
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP091422

Quant Time: Sep 23 01:51:31 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 20 09:45:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	71902	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	118255	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	110585	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	54725	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	71612	54.299	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	108.600%
35) Dibromofluoromethane	5.391	113	61225	50.092	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.180%
50) Toluene-d8	8.647	98	238036	50.523	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.040%
62) 4-Bromofluorobenzene	11.079	95	76358	47.847	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.700%

Target Compounds					Qvalue	
39) Methylcyclohexane	7.379	83	2473	1.069	ug/l	# 85
61) 1,2-Dibromoethane	9.616	107	2774	1.709	ug/l	93
67) Ethyl Benzene	10.195	91	1589	0.264	ug/l	97
68) m/p-Xylenes	10.299	106	11069	4.575	ug/l	95
69) o-Xylene	10.640	106	42411	18.363	ug/l	96
78) n-propylbenzene	11.299	91	1416	0.301	ug/l	# 52
80) 1,3,5-Trimethylbenzene	11.451	105	36693	10.580	ug/l	100
84) 1,2,4-Trimethylbenzene	11.750	105	37497	10.978	ug/l	100
86) p-Isopropyltoluene	12.012	119	2244	0.625	ug/l	84
95) Naphthalene	13.774	128	20599	5.251	ug/l	# 97

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

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