

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062822\
 Data File : VX029844.D
 Acq On : 28 Jun 2022 21:22
 Operator : JC/MD
 Sample : N3494-29
 Misc : 5.60g/5.0mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :

Manual Integrations
 APPROVED

Reviewed By : Mahesh Dadoda 06/30/2022
 Supervised By : Semsettin Yesilyurt 06/30/2022

Quant Time: Jun 29 06:09:15 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 05:54:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	273006	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	462243	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	387653	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	180227	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	147625	42.718	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	85.440%		
35) Dibromofluoromethane	5.391	113	123984	42.919	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	85.840%		
50) Toluene-d8	8.653	98	521323	48.105	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	96.220%		
62) 4-Bromofluorobenzene	11.085	95	176743	46.138	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	92.280%		
Target Compounds					Qvalue	
18) Methyl Acetate	2.727	43	7552	1.449	ug/l	100
40) Benzene	6.031	78	62157	5.003	ug/l	96
65) Chlorobenzene	10.086	112	3629781	466.455	ug/l	95
67) Ethyl Benzene	10.195	91	28360	2.095	ug/l	98
68) m/p-Xylenes	10.305	106	24993	4.815	ug/l	99
69) o-Xylene	10.646	106	6742	1.282	ug/l	94
87) 1,3-Dichlorobenzene	11.975	146	9999	1.660	ug/l	90
88) 1,4-Dichlorobenzene	12.049	146	41023	6.676	ug/l	96
91) 1,2-Dichlorobenzene	12.341	146	100508	16.988	ug/l	98
95) Naphthalene	13.780	128	19588	1.346	ug/l #	91

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

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