

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040423\
 Data File : BG056976.D
 Acq On : 4 Apr 2023 22:23
 Operator : CG/JU
 Sample : 02084-03DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-6ADL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/05/2023
 Supervised By : Jagrut Upadhyay 04/05/2023

Quant Time: Apr 05 04:12:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 03:23:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.413	152	13250	20.000	ng	0.00
21) Naphthalene-d8	11.256	136	49310	20.000	ng	# 0.00
39) Acenaphthene-d10	15.028	164	41073	20.000	ng	0.00
64) Phenanthrene-d10	17.765	188	116195	20.000	ng	# 0.00
76) Chrysene-d12	22.083	240	119803	20.000	ng	-0.01
86) Perylene-d12	25.619	264	125725	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.922	112	7739	9.348	ng	0.00
7) Phenol-d6	7.538	99	7016	5.666	ng	0.00
23) Nitrobenzene-d5	9.588	82	18684	15.504	ng	0.00
42) 2,4,6-Tribromophenol	16.502	330	17979	27.401	ng	0.00
45) 2-Fluorobiphenyl	13.653	172	49476	16.334	ng	0.00
79) Terphenyl-d14	20.332	244	113162	15.984	ng	0.00
Target Compounds						Qvalue
17) 2-Methylphenol	8.836	107	16628	18.996	ng	89
20) 3+4-Methylphenols	9.159	107	4271	3.489	ng	93
27) 2,4-Dimethylphenol	10.416	122	4500	5.359	ng	# 69
31) Naphthalene	11.303	128	54207	20.168	ng	99
37) 2-Methylnaphthalene	12.878	142	11559m	5.419	ng	
38) 1-Methylnaphthalene	13.095	142	6863m	3.426	ng	

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

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