

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050322\
 Data File : BG053417.D
 Acq On : 3 May 2022 14:56
 Operator : CG/JU
 Sample : N2624-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 N47992

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/04/2022
 Supervised By : Jagrut Upadhyay 05/04/2022

Quant Time: May 04 06:06:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 04 06:03:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.102	152	19933	20.000	ng	0.00
21) Naphthalene-d8	10.910	136	80255	20.000	ng	# 0.00
39) Acenaphthene-d10	14.722	164	63785	20.000	ng	0.00
64) Phenanthrene-d10	17.466	188	167609	20.000	ng	0.00
76) Chrysene-d12	21.748	240	177937	20.000	ng	0.00
86) Perylene-d12	25.044	264	182485	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.670	112	86648	74.515	ng	0.00
7) Phenol-d6	7.274	99	80003	44.615	ng	0.01
23) Nitrobenzene-d5	9.271	82	214688	108.617	ng	0.00
42) 2,4,6-Tribromophenol	16.214	330	167544	165.102	ng	0.00
45) 2-Fluorobiphenyl	13.348	172	487114	105.304	ng	0.00
79) Terphenyl-d14	20.074	244	888423	84.325	ng	0.00
Target Compounds						Qvalue
9) Benzaldehyde	7.238	77	9254	8.633	ng	95
15) Benzyl Alcohol	8.448	79	6952063	4361.683	ng	95
32) Benzoic acid	10.199	122	10715m	16.848	ng	
52) Acenaphthene	14.787	154	9546	2.454	ng	97

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

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