

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050322\
 Data File : BG053422.D
 Acq On : 3 May 2022 18:09
 Operator : CG/JU
 Sample : N2641-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-24

Manual Integrations
 APPROVED

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

Quant Time: May 04 06:07:24 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	8.101	152	17651	20.000	ng	# 0.00
21) Naphthalene-d8	10.909	136	65248	20.000	ng	# 0.00
39) Acenaphthene-d10	14.722	164	50661	20.000	ng	0.00
64) Phenanthrene-d10	17.471	188	146409	20.000	ng	0.00
76) Chrysene-d12	21.747	240	162312	20.000	ng	0.00
86) Perylene-d12	25.037	264	163303	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.669	112	182478	177.214	ng	0.00
7) Phenol-d6	7.261	99	241076	151.820	ng	0.00
23) Nitrobenzene-d5	9.264	82	199732	124.292	ng	0.00
42) 2,4,6-Tribromophenol	16.214	330	166603	206.704	ng	0.00
45) 2-Fluorobiphenyl	13.347	172	436493	118.806	ng	0.00
79) Terphenyl-d14	20.073	244	1055290	109.806	ng	0.00
Target Compounds						
77) Benzidine	19.703	184	27162m	5.865	ng	Qvalue

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

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