

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051922\
 Data File : BG053636.D
 Acq On : 19 May 2022 21:31
 Operator : CG/JU
 Sample : N2871-16 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P017-SS004-0006-01

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/20/2022
 Supervised By : Jagrut Upadhyay 05/20/2022

Quant Time: May 20 03:58:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 20 03:55:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.082	152	36620	20.000	ng	0.00
21) Naphthalene-d8	10.890	136	143192	20.000	ng	0.00
39) Acenaphthene-d10	14.702	164	103110	20.000	ng	0.00
64) Phenanthrene-d10	17.452	188	210033	20.000	ng	0.00
76) Chrysene-d12	21.728	240	179432	20.000	ng	0.00
86) Perylene-d12	25.012	264	182438	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.656	112	13334	6.180	ng	0.01
7) Phenol-d6	7.254	99	19225m	5.871	ng	0.01
23) Nitrobenzene-d5	9.245	82	13401	3.984	ng	0.00
42) 2,4,6-Tribromophenol	16.206	330	5756	3.785	ng	0.01
45) 2-Fluorobiphenyl	13.328	172	32297	4.577	ng	0.00
79) Terphenyl-d14	20.054	244	47466	4.927	ng	0.00

Target Compounds Qvalue

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

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