

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062722\
 Data File : BG054100.D
 Acq On : 27 Jun 2022 20:26
 Operator : CG/JU
 Sample : N3494-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP-EXC-1-4

Quant Time: Jun 28 04:58:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 00:27:54 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 06/28/2022
 Supervised By :Jagrut Upadhyay 06/28/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.070	152	24923	20.000	ng	0.00
21) Naphthalene-d8	10.878	136	94922	20.000	ng	0.00
39) Acenaphthene-d10	14.697	164	69322	20.000	ng	0.00
64) Phenanthrene-d10	17.441	188	185245	20.000	ng	0.00
76) Chrysene-d12	21.713	240	229587	20.000	ng	0.00
86) Perylene-d12	24.968	264	224560	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.643	112	190347	146.020	ng	0.00
7) Phenol-d6	7.230	99	229628	129.852	ng	0.00
23) Nitrobenzene-d5	9.233	82	178161	105.476	ng	0.00
42) 2,4,6-Tribromophenol	16.184	330	163255	146.577	ng	0.00
45) 2-Fluorobiphenyl	13.322	172	451624	94.735	ng	0.00
79) Terphenyl-d14	20.050	244	1148573	100.760	ng	0.00
Target Compounds						
32) Benzoic acid	10.203	122	3704m	16.754	ng	Qvalue

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

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