

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080422\
 Data File : BG054527.D
 Acq On : 4 Aug 2022 20:37
 Operator : CG/JU
 Sample : N3930-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-99

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/05/2022
 Supervised By : Jagrut Upadhyay 08/05/2022

Quant Time: Aug 05 05:03:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.982	152	21503	20.00	ng	0.00
21) Naphthalene-d8	10.791	136	81809	20.00	ng	# 0.00
39) Acenaphthene-d10	14.622	164	59523	20.00	ng	0.00
64) Phenanthrene-d10	17.371	188	137929	20.00	ng	# 0.00
76) Chrysene-d12	21.637	240	144913	20.00	ng	# 0.00
86) Perylene-d12	24.851	264	123203	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.574	112	62161	60.36	ng	0.01
7) Phenol-d6	7.183	99	82339	58.35	ng	0.01
23) Nitrobenzene-d5	9.146	82	59086	39.33	ng	0.00
42) 2,4,6-Tribromophenol	16.120	330	67661	54.14	ng	0.00
45) 2-Fluorobiphenyl	13.241	172	176548	41.84	ng	0.00
79) Terphenyl-d14	19.974	244	313399	42.90	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	19.422	202	25800	2.76	ng	95
78) Pyrene	19.786	202	30650	3.69	ng	95
88) Benzo(b)fluoranthene	23.834	252	20941	2.87	ng	# 73
90) Benzo(a)pyrene	24.704	252	14519	2.33	ng	# 90

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

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