

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
 Data File : BG051735.D
 Acq On : 22 Dec 2021 4:23
 Operator : CG/JU
 Sample : PB141474BL
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK474

Quant Time: Dec 22 06:08:59 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 14 14:19:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.270	152	42290	20.000	ng/u1	-0.02
20) Naphthalene-d8	11.107	136	175903	20.000	ng/u1	#-0.02
38) Acenaphthene-d10	14.908	164	126507	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.657	188	295371	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.975	240	289058	20.000	ng/u1	#-0.01
88) Perylene-d12	25.476	264	274469	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.600	96	4633	4.550	ng/uL	0.00
4) Pyridine-d5	4.028	84	62443	20.295	ng/u1	-0.02
7) Phenol-d5	7.406	99	100620	26.648	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.576	67	58493	26.058	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.800	132	69616	26.509	ng/u1	-0.01
15) 4-Methylphenol-d8	8.969	113	78025	26.733	ng/u1	-0.01
21) Nitrobenzene-d5	9.439	128	35745	27.717	ng/u1	-0.02
24) 2-Nitrophenol-d4	10.173	143	38653	27.417	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.719	165	78616	25.715	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.230	131	42406	10.535	ng/u1	-0.01
46) Dimethylphthalate-d6	14.297	166	257942	25.474	ng/u1	-0.02
49) Acenaphthylene-d8	14.602	160	320707	25.503	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.072	143	35953	21.835	ng/u1	0.00
60) Fluorene-d10	15.895	176	223079	24.590	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.995	200	33821	21.977	ng/u1	-0.02
73) Anthracene-d10	17.757	188	377466	26.790	ng/u1	0.00
81) Pyrene-d10	20.030	212	454305	26.114	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.235	264	420017	29.084	ng/u1	-0.02

Target Compounds Qvalue

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

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