

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050836.D
 Acq On : 3 Nov 2021 5:19
 Operator : CG/JU
 Sample : PB140350BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK350

Quant Time: Nov 03 08:05:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 02 14:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.232	152	41415	20.000	ng/u1	0.00
20) Naphthalene-d8	11.052	136	207467	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.854	164	157323	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.598	188	373742	20.000	ng/u1	0.00
79) Chrysene-d12	21.893	240	342751	20.000	ng/u1	#-0.01
88) Perylene-d12	25.300	264	328749	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.585	96	5268	4.105	ng/uL	0.00
4) Pyridine-d5	4.014	84	73390	19.118	ng/u1	0.00
7) Phenol-d5	7.369	99	86637	19.609	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.539	67	59695	20.916	ng/u1	0.00
11) 2-Chlorophenol-d4	7.756	132	61968	20.238	ng/u1	0.00
15) 4-Methylphenol-d8	8.925	113	68781	19.774	ng/u1	0.00
21) Nitrobenzene-d5	9.401	128	34549	19.596	ng/u1	0.00
24) 2-Nitrophenol-d4	10.130	143	38999	19.894	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.671	165	58904	17.837	ng/u1	0.00
31) 4-Chloroaniline-d4	11.188	131	95101	19.017	ng/u1	0.00
46) Dimethylphthalate-d6	14.249	166	250465	20.809	ng/u1	0.00
49) Acenaphthylene-d8	14.548	160	288917	19.267	ng/u1	0.00
54) 4-Nitrophenol-d4	15.036	143	36989	16.949	ng/u1	0.00
60) Fluorene-d10	15.841	176	209080	19.609	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	37669	16.624	ng/u1	0.00
73) Anthracene-d10	17.698	188	354329	20.052	ng/u1	0.00
81) Pyrene-d10	19.971	212	416537	18.816	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.060	264	343375	18.894	ng/u1	-0.01

Target Compounds Qvalue

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

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