

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050992.D
 Acq On : 12 Nov 2021 8:50
 Operator : CG/JU
 Sample : PB140665BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK665

Quant Time: Nov 12 11:48:49 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 11 12:40:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.229	152	45893	20.000	ng/ul	0.00
20) Naphthalene-d8	11.055	136	215017	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.851	164	148845	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.594	188	344364	20.000	ng/ul	0.00
79) Chrysene-d12	21.889	240	314494	20.000	ng/ul	0.00
88) Perylene-d12	25.291	264	306083	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.587	96	8007	5.631	ng/uL	0.00
4) Pyridine-d5	4.010	84	116458	27.377	ng/ul	0.00
7) Phenol-d5	7.371	99	133289	27.224	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	89643	28.344	ng/ul	0.00
11) 2-Chlorophenol-d4	7.753	132	95232	28.067	ng/ul	0.00
15) 4-Methylphenol-d8	8.922	113	104643	27.149	ng/ul	0.00
21) Nitrobenzene-d5	9.404	128	52162	28.547	ng/ul	0.00
24) 2-Nitrophenol-d4	10.127	143	58820	28.951	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.667	165	91523	26.742	ng/ul	0.00
31) 4-Chloroaniline-d4	11.184	131	143156	27.621	ng/ul	0.00
46) Dimethylphthalate-d6	14.245	166	349447	30.687	ng/ul	0.00
49) Acenaphthylene-d8	14.551	160	412654	29.086	ng/ul	0.00
54) 4-Nitrophenol-d4	15.039	143	60149	29.131	ng/ul	0.00
60) Fluorene-d10	15.838	176	293740	29.118	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.955	200	53653	25.698	ng/ul	0.00
73) Anthracene-d10	17.694	188	490684	30.138	ng/ul	0.00
81) Pyrene-d10	19.968	212	578949	28.502	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.056	264	486149	28.731	ng/ul	0.00

Target Compounds Qvalue

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

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