

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050865.D
 Acq On : 4 Nov 2021 3:09
 Operator : CG/JU
 Sample : PB140458BL
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK458

Quant Time: Nov 07 08:07:46 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	8.229	152	28708	20.000	ng/u1	0.00
20) Naphthalene-d8	11.055	136	131825	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.850	164	92525	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.594	188	215173	20.000	ng/u1	-0.01
79) Chrysene-d12	21.889	240	190843	20.000	ng/u1	#-0.02
88) Perylene-d12	25.285	264	187095	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.581	96	5329	5.991	ng/uL	0.00
4) Pyridine-d5	4.010	84	80567	30.278	ng/u1	0.00
7) Phenol-d5	7.371	99	92357	30.156	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	61646	31.160	ng/u1	0.00
11) 2-Chlorophenol-d4	7.753	132	65305	30.768	ng/u1	-0.01
15) 4-Methylphenol-d8	8.922	113	71984	29.856	ng/u1	-0.01
21) Nitrobenzene-d5	9.398	128	36465	32.551	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.126	143	39086	31.378	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.667	165	61106	29.122	ng/u1	0.00
31) 4-Chloroaniline-d4	11.184	131	100486	31.624	ng/u1	0.00
46) Dimethylphthalate-d6	14.245	166	238564	33.701	ng/u1	0.00
49) Acenaphthylene-d8	14.551	160	286838	32.524	ng/u1	0.00
54) 4-Nitrophenol-d4	15.038	143	36163	28.176	ng/u1	0.00
60) Fluorene-d10	15.837	176	203297	32.420	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.955	200	33435	25.630	ng/u1	0.00
73) Anthracene-d10	17.694	188	340344	33.455	ng/u1	-0.01
81) Pyrene-d10	19.974	212	396943	32.203	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.056	264	331622	32.063	ng/u1	-0.02

Target Compounds Qvalue

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

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