

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
 Data File : BG050205.D
 Acq On : 8 Sep 2021 19:22
 Operator : CG/JU
 Sample : PB138916BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK916

Quant Time: Sep 09 00:56:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 07 18:32:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	49668	20.000	ng/ul	0.00
20) Naphthalene-d8	10.762	136	149811	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.610	164	100985	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	223653	20.000	ng/ul	0.00
79) Chrysene-d12	21.636	240	240003	20.000	ng/ul	0.00
88) Perylene-d12	24.849	264	235417	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.308	96	6650	6.619	ng/uL	0.00
4) Pyridine-d5	3.737	84	127382	35.719	ng/ul	0.00
7) Phenol-d5	7.120	99	118058	30.614	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.267	67	91321	40.547	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.479	132	94293	29.462	ng/ul	0.00
15) 4-Methylphenol-d8	8.671	113	101735	33.553	ng/ul	0.00
21) Nitrobenzene-d5	9.123	128	36849	35.835	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	43167	34.523	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	75324	31.563	ng/ul	0.00
31) 4-Chloroaniline-d4	10.909	131	97617	33.201	ng/ul	0.00
46) Dimethylphthalate-d6	14.011	166	304331	39.911	ng/ul	0.00
49) Acenaphthylene-d8	14.299	160	306428	33.750	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.845	143	27092	29.446	ng/ul	0.00
60) Fluorene-d10	15.603	176	206085	32.550	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.744	200	32793	24.307	ng/ul	0.00
73) Anthracene-d10	17.459	188	327777	32.647	ng/ul	0.00
81) Pyrene-d10	19.756	212	399677	33.433	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.626	264	392426	31.489	ng/ul	-0.01

Target Compounds	Qvalue

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

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