

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092821\
 Data File : BG050287.D
 Acq On : 28 Sep 2021 14:47
 Operator : CG/JU
 Sample : PB139009BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK009

Quant Time: Sep 28 15:26:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG092521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 27 03:10:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.288	152	35528	20.000	ng/u1	0.00
20) Naphthalene-d8	11.114	136	139408	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.904	164	111280	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.648	188	404322	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.949	240	449816	20.000	ng/u1	# 0.00
88) Perylene-d12	25.392	264	477333	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.652	96	5997	8.163	ng/uL	0.00
4) Pyridine-d5	4.087	84	81624	36.863	ng/u1	0.00
7) Phenol-d5	7.419	99	91893	32.576	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.601	67	57350	35.028	ng/u1	0.00
11) 2-Chlorophenol-d4	7.812	132	65875	32.129	ng/u1	0.00
15) 4-Methylphenol-d8	8.976	113	69495	31.273	ng/u1	0.00
21) Nitrobenzene-d5	9.457	128	32561	37.808	ng/u1	0.00
24) 2-Nitrophenol-d4	10.180	143	33560	35.281	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.721	165	63301	28.527	ng/u1	0.00
31) 4-Chloroaniline-d4	11.238	131	94514	32.423	ng/u1	0.00
46) Dimethylphthalate-d6	14.299	166	404851	53.530	ng/u1	0.00
49) Acenaphthylene-d8	14.598	160	313049	32.928	ng/u1	0.00
54) 4-Nitrophenol-d4	15.063	143	63192	49.521	ng/u1	0.00
60) Fluorene-d10	15.891	176	337894	48.797	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.991	200	68163	39.793	ng/u1	0.00
73) Anthracene-d10	17.742	188	605829	35.179	ng/u1	0.00
81) Pyrene-d10	20.016	212	745686	30.619	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.157	264	811184	33.332	ng/u1	0.00

Target Compounds	Qvalue

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

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