

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092821\
 Data File : BG050285.D
 Acq On : 28 Sep 2021 13:20
 Operator : CG/JU
 Sample : PB139008BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK008

Quant Time: Sep 28 13:59:04 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	35807	20.000	ng/ul	0.00
20) Naphthalene-d8	11.114	136	136376	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.904	164	96653	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.648	188	222990	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.949	240	250586	20.000	ng/ul	# 0.00
88) Perylene-d12	25.392	264	256805	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.652	96	5494	7.420	ng/uL	0.00
4) Pyridine-d5	4.081	84	80479	36.063	ng/ul	-0.01
7) Phenol-d5	7.413	99	92599	32.571	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.601	67	57603	34.908	ng/ul	0.00
11) 2-Chlorophenol-d4	7.812	132	65832	31.858	ng/ul	0.00
15) 4-Methylphenol-d8	8.976	113	68106	30.409	ng/ul	0.00
21) Nitrobenzene-d5	9.457	128	31681	37.604	ng/ul	0.00
24) 2-Nitrophenol-d4	10.180	143	33839	36.365	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.721	165	64176	29.565	ng/ul	0.00
31) 4-Chloroaniline-d4	11.238	131	91327	32.026	ng/ul	0.00
46) Dimethylphthalate-d6	14.299	166	208725	31.774	ng/ul	0.00
49) Acenaphthylene-d8	14.598	160	255597	30.954	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	34399	31.037	ng/ul	0.00
60) Fluorene-d10	15.891	176	186184	30.957	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.991	200	30108	31.870	ng/ul	0.00
73) Anthracene-d10	17.748	188	310702	32.713	ng/ul	0.00
81) Pyrene-d10	20.016	212	391804	28.879	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.157	264	409567	31.281	ng/ul	0.00

Target Compounds Qvalue

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

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