

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093021\
 Data File : BG050394.D
 Acq On : 2 Oct 2021 18:19
 Operator : CG/JU
 Sample : PB139405BL
 Misc :
 ALS Vial : 80 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK405

Quant Time: Oct 04 01:08:06 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG093021.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 30 16:27:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.273	152	43792	20.00	ng/ul	0.00
20) Naphthalene-d8	11.093	136	185428	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.889	164	124104	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.627	188	269247	20.00	ng/ul	0.00
79) Chrysene-d12	21.928	240	252274	20.00	ng/ul	-0.02
88) Perylene-d12	25.353	264	249695	20.00	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.637	96	8244	6.92	ng/uL	0.00
4) Pyridine-d5	4.066	84	120893	34.04	ng/ul	0.00
7) Phenol-d5	7.398	99	134495	35.15	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.586	67	90629	37.64	ng/ul	0.00
11) 2-Chlorophenol-d4	7.797	132	95283	35.54	ng/ul	0.00
15) 4-Methylphenol-d8	8.955	113	104980	36.25	ng/ul	0.00
21) Nitrobenzene-d5	9.436	128	51247	39.51	ng/ul	0.00
24) 2-Nitrophenol-d4	10.165	143	54065	38.83	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.700	165	92310	33.58	ng/ul	0.00
31) 4-Chloroaniline-d4	11.222	131	139006	35.39	ng/ul	0.00
46) Dimethylphthalate-d6	14.278	166	302605	35.63	ng/ul	-0.01
49) Acenaphthylene-d8	14.583	160	377460	35.39	ng/ul	0.00
54) 4-Nitrophenol-d4	15.047	143	55953	34.20	ng/ul	0.00
60) Fluorene-d10	15.876	176	266470	34.82	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	41935	31.23	ng/ul	0.00
73) Anthracene-d10	17.727	188	423813	37.26	ng/ul	0.00
81) Pyrene-d10	20.000	212	491353	36.63	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.118	264	439782	35.03	ng/ul	-0.02

Target Compounds Qvalue

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

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