

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG103123\
 Data File : BG059556.D
 Acq On : 31 Oct 2023 10:10
 Operator : MA/JU
 Sample : PB156466BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK466

Quant Time: Nov 01 03:30:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 28 03:56:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.359	152	77797	20.000	ng/u1	0.00
20) Naphthalene-d8	11.209	136	446696	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.987	164	267738	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.736	188	609034	20.000	ng/u1	0.00
79) Chrysene-d12	22.073	240	513821	20.000	ng/u1	0.00
88) Perylene-d12	25.686	264	553528	20.000	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.659	96	10753	5.171	ng/uL	0.00
4) Pyridine-d5	4.088	84	147990	25.469	ng/u1	0.00
7) Phenol-d5	7.466	99	263386	32.946	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.672	67	95236	25.936	ng/u1	0.00
11) 2-Chlorophenol-d4	7.872	132	170716	26.322	ng/u1	0.00
15) 4-Methylphenol-d8	9.041	113	175316	27.641	ng/u1	0.00
21) Nitrobenzene-d5	9.558	128	106135	33.416	ng/u1	0.00
24) 2-Nitrophenol-d4	10.275	143	123644	37.920	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.798	165	216789	30.177	ng/u1	0.00
31) 4-Chloroaniline-d4	11.344	131	328996	27.820	ng/u1	0.00
46) Dimethylphthalate-d6	14.382	166	697547	28.757	ng/u1	0.00
49) Acenaphthylene-d8	14.687	160	818827	28.765	ng/u1	0.00
54) 4-Nitrophenol-d4	15.151	143	131202	31.542	ng/u1	0.00
60) Fluorene-d10	15.974	176	596338	28.767	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.074	200	82160	28.486	ng/u1	0.00
73) Anthracene-d10	17.836	188	948352	28.742	ng/u1	0.00
81) Pyrene-d10	20.110	212	1055958	29.566	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.433	264	998803	30.643	ng/u1	0.00

Target Compounds Qvalue

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

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