

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102721\
 Data File : BG050739.D
 Acq On : 27 Oct 2021 10:53
 Operator : CG/JU
 Sample : PB140133BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK133

Quant Time: Oct 27 11:32:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 26 18:31:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.247	152	32767	20.000	ng/u1	0.00
20) Naphthalene-d8	11.073	136	150498	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.874	164	100859	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.624	188	222752	20.000	ng/u1	0.00
79) Chrysene-d12	21.931	240	184000	20.000	ng/u1	# 0.00
88) Perylene-d12	25.356	264	178231	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.605	96	5276	6.344	ng/uL	0.00
4) Pyridine-d5	4.028	84	74558	28.311	ng/u1	0.00
7) Phenol-d5	7.389	99	89244	29.309	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.559	67	59269	30.254	ng/u1	0.00
11) 2-Chlorophenol-d4	7.777	132	63544	29.710	ng/u1	0.00
15) 4-Methylphenol-d8	8.946	113	70396	29.943	ng/u1	0.00
21) Nitrobenzene-d5	9.422	128	33820	29.441	ng/u1	0.00
24) 2-Nitrophenol-d4	10.145	143	38301	29.567	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.685	165	57914	25.457	ng/u1	0.00
31) 4-Chloroaniline-d4	11.208	131	92151	27.600	ng/u1	0.00
46) Dimethylphthalate-d6	14.269	166	230545	32.366	ng/u1	0.00
49) Acenaphthylene-d8	14.569	160	259617	29.348	ng/u1	0.00
54) 4-Nitrophenol-d4	15.056	143	36643	28.425	ng/u1	0.00
60) Fluorene-d10	15.861	176	188603	30.811	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.979	200	31740	24.349	ng/u1	0.00
73) Anthracene-d10	17.724	188	317981	33.294	ng/u1	0.00
81) Pyrene-d10	19.998	212	357175	34.063	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.115	264	288002	31.407	ng/u1	0.00

Target Compounds	Qvalue

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

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