

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081121\
 Data File : BG049795.D
 Acq On : 11 Aug 2021 19:39
 Operator : CG/JU
 Sample : PB138195BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK195

Quant Time: Aug 12 01:31:45 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 10 16:55:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.036	152	25336	20.000	ng/ul	0.00
20) Naphthalene-d8	10.856	136	105610	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.686	164	75052	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.441	188	167166	20.000	ng/ul	0.00
79) Chrysene-d12	21.718	240	156375	20.000	ng/ul	0.00
88) Perylene-d12	24.995	264	153101	20.000	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.425	96	3037	5.604	ng/uL	0.01
4) Pyridine-d5	3.848	84	43380	26.676	ng/ul	0.00
7) Phenol-d5	7.196	99	66793	29.048	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.355	67	37391	28.310	ng/ul	0.00
11) 2-Chlorophenol-d4	7.566	132	51382	31.221	ng/ul	0.00
15) 4-Methylphenol-d8	8.753	113	53170	30.327	ng/ul	0.00
21) Nitrobenzene-d5	9.205	128	26460	31.806	ng/ul	0.00
24) 2-Nitrophenol-d4	9.928	143	34101	35.893	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.480	165	52557	29.338	ng/ul	0.00
31) 4-Chloroaniline-d4	10.991	131	75027	30.593	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	210043	35.663	ng/ul	0.00
49) Acenaphthylene-d8	14.380	160	232810	33.225	ng/ul	0.00
54) 4-Nitrophenol-d4	14.892	143	29982	31.485	ng/ul	0.00
60) Fluorene-d10	15.679	176	171631	33.901	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.802	200	33392	29.679	ng/ul	0.00
73) Anthracene-d10	17.541	188	283879	35.976	ng/ul	0.00
81) Pyrene-d10	19.826	212	316662	36.753	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.766	264	292869	34.911	ng/ul	0.00

Target Compounds	Qvalue

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

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