

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050147.D
 Acq On : 4 Sep 2021 2:29
 Operator : CG/JU
 Sample : PB138916BL
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK916

Quant Time: Sep 04 03:23:35 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 14:59:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	34688	20.000	ng/u1	0.00
20) Naphthalene-d8	10.766	136	138577	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.608	164	94340	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.363	188	204341	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.634	240	204330	20.000	ng/u1	-0.01
88) Perylene-d12	24.853	264	199952	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	4532	6.737	ng/uL	0.00
4) Pyridine-d5	3.734	84	56803	28.034	ng/u1	0.00
7) Phenol-d5	7.124	99	87066	29.558	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.271	67	49959	30.529	ng/u1	0.00
11) 2-Chlorophenol-d4	7.482	132	66892	30.350	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	70242	30.279	ng/u1	0.00
21) Nitrobenzene-d5	9.121	128	34834	32.423	ng/u1	0.00
24) 2-Nitrophenol-d4	9.850	143	40652	31.777	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.402	165	66811	29.281	ng/u1	0.00
31) 4-Chloroaniline-d4	10.907	131	89601	28.851	ng/u1	-0.01
46) Dimethylphthalate-d6	14.015	166	226329	31.348	ng/u1	0.00
49) Acenaphthylene-d8	14.302	160	267360	31.583	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.849	143	25985	24.876	ng/u1	0.00
60) Fluorene-d10	15.606	176	182946	29.886	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.747	200	27903	21.225	ng/u1	0.00
73) Anthracene-d10	17.463	188	297046	32.372	ng/u1	-0.01
81) Pyrene-d10	19.754	212	346197	31.353	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.635	264	328784	30.994	ng/u1	-0.01

Target Compounds	Qvalue

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

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