

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
 Data File : BG048210.D
 Acq On : 19 Feb 2021 19:28
 Operator : CG/JU
 Sample : PB134690BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK90

Quant Time: Feb 19 22:10:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 19 11:17:07 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	37944	20.00	ng/ul	0.00
18) Naphthalene-d8	10.92	136	152429	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	117528	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.48	188	324274	20.00	ng/ul	0.00
78) Chrysene-d12	21.76	240	358092	20.00	ng/ul	0.00
86) Perylene-d12	25.04	264	352847	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.50	96	6301	6.91	ng/uL	0.00
5) Phenol-d5	7.29	99	129109	40.38	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.43	67	79148	41.56	ng/ul	0.00
9) 2-Chlorophenol-d4	7.65	132	91721	40.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.83	113	105067	41.89	ng/ul	0.00
19) Nitrobenzene-d5	9.28	128	47084	41.35	ng/ul	0.00
22) 2-Nitrophenol-d4	10.00	143	59498	43.20	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	105113	38.97	ng/ul	0.00
29) 4-Chloroaniline-d4	11.07	131	133056	49.92	ng/ul	0.00
44) Dimethylphthalate-d6	14.14	166	396532	45.55	ng/ul	0.00
47) Acenaphthylene-d8	14.43	160	452578	42.53	ng/ul	0.00
52) 4-Nitrophenol-d4	14.97	143	64739	43.77	ng/ul	0.00
58) Fluorene-d10	15.72	176	341446	43.97	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.85	200	81023	40.75	ng/ul	0.00
71) Anthracene-d10	17.58	188	627030	45.10	ng/ul	0.00
79) Pyrene-d10	19.86	212	774294	43.30	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.82	264	799986	44.44	ng/ul	0.00

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

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