

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120420\
 Data File : BG047577.D
 Acq On : 4 Dec 2020 19:17
 Operator : CG/JU
 Sample : PB133222BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK22

Quant Time: Dec 05 01:42:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 05 00:50:33 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	28547	20.00	ng/ul	-0.02
18) Naphthalene-d8	11.08	136	112451	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.86	164	92569	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.60	188	234816	20.00	ng/ul	-0.02
78) Chrysene-d12	21.88	240	217119	20.00	ng/ul	-0.02
86) Perylene-d12	25.27	264	222226	20.00	ng/ul	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.59	96	4063	5.17	ng/uL	-0.03
5) Phenol-d5	7.39	99	87292	30.49	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.56	67	50408	30.87	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.78	132	70860	36.90	ng/ul	-0.02
13) 4-Methylphenol-d8	8.94	113	74810	34.00	ng/ul	-0.02
19) Nitrobenzene-d5	9.41	128	36700	39.51	ng/ul	-0.02
22) 2-Nitrophenol-d4	10.14	143	43662	43.08	ng/ul	-0.03
26) 2,4-Dichlorophenol-d3	10.68	165	84078	36.23	ng/ul	-0.02
29) 4-Chloroaniline-d4	11.20	131	102131	43.52	ng/ul	-0.02
44) Dimethylphthalate-d6	14.26	166	316736	39.86	ng/ul	-0.02
47) Acenaphthylene-d8	14.56	160	337273	36.72	ng/ul	-0.02
52) 4-Nitrophenol-d4	15.03	143	45464	37.50	ng/ul	0.00
58) Fluorene-d10	15.85	176	271105	37.24	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.95	200	59056	44.18	ng/ul	-0.02
71) Anthracene-d10	17.70	188	472260	40.54	ng/ul	-0.02
79) Pyrene-d10	19.97	212	519737	41.14	ng/ul	-0.01
90) Benzo(a)pyrene-d12	25.04	264	506346	39.59	ng/ul	-0.02

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

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