

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\
 Data File : BG047503.D
 Acq On : 1 Dec 2020 19:08
 Operator : CG/JU
 Sample : PB133291BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK91

Manual Integrations
 APPROVED

mohammad
 12/2/2020 2:26:07 PM

Quant Time: Dec 02 07:41:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 23 13:14:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	24663	20.00	ng/ul	-0.01
18) Naphthalene-d8	11.09	136	99149	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.88	164	82097	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.61	188	223772	20.00	ng/ul	0.00
78) Chrysene-d12	21.90	240	233588	20.00	ng/ul	0.00
86) Perylene-d12	25.29	264	227135	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.60	96	3383m	4.99	ng/uL	-0.02
5) Phenol-d5	7.40	99	65559	26.51	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.58	67	41337	29.30	ng/ul	0.00
9) 2-Chlorophenol-d4	7.79	132	48182	29.05	ng/ul	-0.01
13) 4-Methylphenol-d8	8.95	113	57500	30.25	ng/ul	-0.01
19) Nitrobenzene-d5	9.44	128	27716	33.84	ng/ul	0.00
22) 2-Nitrophenol-d4	10.16	143	33788	37.81	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.71	165	64244	31.40	ng/ul	0.00
29) 4-Chloroaniline-d4	11.22	131	84611	40.89	ng/ul	-0.01
44) Dimethylphthalate-d6	14.27	166	301900	42.84	ng/ul	-0.01
47) Acenaphthylene-d8	14.58	160	310651	38.14	ng/ul	0.00
52) 4-Nitrophenol-d4	15.04	143	45259	42.09	ng/ul	-0.01
58) Fluorene-d10	15.86	176	253793	39.31	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.96	200	53708	42.17	ng/ul	-0.01
71) Anthracene-d10	17.71	188	452627m	40.77	ng/ul	0.00
79) Pyrene-d10	19.98	212	542871	39.94	ng/ul	0.00
90) Benzo(a)pyrene-d12	25.06	264	531733	40.67	ng/ul	-0.02

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

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