

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102220\
 Data File : BG047029.D
 Acq On : 22 Oct 2020 13:05
 Operator : CG/JU
 Sample : PB132535BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK35

Quant Time: Oct 22 13:45:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 22 05:48:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	46184	20.00	ng/ul	0.00
18) Naphthalene-d8	10.87	136	187400	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	136649	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	329945	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	370757	20.00	ng/ul	0.00
86) Perylene-d12	24.93	264	313176	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.47	96	9396	8.21	ng/uL	0.00
5) Phenol-d5	7.21	99	162724	40.60	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	98056	42.53	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	127863	41.02	ng/ul	0.00
13) 4-Methylphenol-d8	8.76	113	130438	40.79	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	64611	41.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	75631	41.22	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	134723	37.63	ng/ul	0.00
29) 4-Chloroaniline-d4	11.00	131	140108	45.89	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	474003	42.19	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	553904	41.67	ng/ul	0.00
52) 4-Nitrophenol-d4	14.87	143	67482	38.15	ng/ul	0.00
58) Fluorene-d10	15.68	176	413495	39.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	86145	37.86	ng/ul	0.00
71) Anthracene-d10	17.53	188	677684	42.38	ng/ul	0.00
79) Pyrene-d10	19.81	212	870089	41.14	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.71	264	726590	42.20	ng/ul	0.00

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

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