

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122320\
 Data File : BG047811.D
 Acq On : 23 Dec 2020 23:18
 Operator : CG/JU
 Sample : L5112-17
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A54T2

Quant Time: Dec 24 00:55:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG120720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 23 13:42:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.20	152	27592	20.00	ng/ul	0.00
18) Naphthalene-d8	11.02	136	100608	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.82	164	75750	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.56	188	194624	20.00	ng/ul	0.00
78) Chrysene-d12	21.82	240	198365	20.00	ng/ul	0.00
86) Perylene-d12	25.18	264	200760	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.53	96	2107	4.07	ng/uL	0.00
5) Phenol-d5	7.33	99	37373	15.83	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.51	67	19563	16.11	ng/ul	0.00
9) 2-Chlorophenol-d4	7.72	132	26527	14.81	ng/ul	0.00
13) 4-Methylphenol-d8	8.89	113	20078	10.71	ng/ul	0.00
19) Nitrobenzene-d5	9.37	128	13924	16.97	ng/ul	0.00
22) 2-Nitrophenol-d4	10.10	143	16806	17.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.63	165	32790	15.27	ng/ul	0.00
29) 4-Chloroaniline-d4	11.15	131	29076	14.75	ng/ul	0.00
44) Dimethylphthalate-d6	14.21	166	105164	15.98	ng/ul	0.00
47) Acenaphthylene-d8	14.51	160	130850	17.41	ng/ul	0.00
52) 4-Nitrophenol-d4	14.98	143	13368	13.70	ng/ul	0.00
58) Fluorene-d10	15.81	176	101067	16.69	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.91	200	17695	13.43	ng/ul	0.00
71) Anthracene-d10	17.65	188	171977	17.72	ng/ul	0.00
79) Pyrene-d10	19.92	212	208207	17.76	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.94	264	210088	17.92	ng/ul	-0.01
Target Compounds						
6) Phenol	7.36	94	3690	1.654	ng/ul	88
45) Dimethylphthalate	14.25	163	12829	1.925	ng/ul#	81

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

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