

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112320\
 Data File : BG047431.D
 Acq On : 24 Nov 2020 3:47
 Operator : CG/JU
 Sample : L4749-22
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0BG1

Manual Integrations
 APPROVED

mohammad
 11/25/2020 12:47:15 PM

Quant Time: Nov 24 06:50:56 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	8.28	152	36113	20.00	ng/ul	0.00
18) Naphthalene-d8	11.11	136	133269	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.88	164	99697	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.62	188	227486m	20.00	ng/ul	0.00
78) Chrysene-d12	21.90	240	203127	20.00	ng/ul	0.00
86) Perylene-d12	25.31	264	211216	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.61	96	72	0.07	ng/uL	0.00
5) Phenol-d5	7.41	99	7520	2.08	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.58	67	4640	2.25	ng/ul	0.00
9) 2-Chlorophenol-d4	7.80	132	6776	2.79	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	5094	1.83	ng/ul	0.00
19) Nitrobenzene-d5	9.45	128	3201m	2.91	ng/ul	0.00
22) 2-Nitrophenol-d4	10.17	143	3707	3.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.71	165	7777	2.83	ng/ul	0.00
29) 4-Chloroaniline-d4	11.22	131	8666	3.12	ng/ul	0.00
44) Dimethylphthalate-d6	14.27	166	23587	2.76	ng/ul	-0.01
47) Acenaphthylene-d8	14.58	160	29346	2.97	ng/ul	0.00
52) 4-Nitrophenol-d4	15.04	143	2637	2.02	ng/ul	-0.01
58) Fluorene-d10	15.87	176	22241	2.84	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.99	200	100	0.08	ng/ul	0.02
71) Anthracene-d10	17.72	188	35512	3.15	ng/ul	0.00
79) Pyrene-d10	19.98	212	41069	3.47	ng/ul	0.00
90) Benzo(a)pyrene-d12	25.07	264	39482	3.25	ng/ul	-0.01

Target Compounds Ovalue

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

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