

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
 Data File : BG046909.D
 Acq On : 15 Oct 2020 1:51
 Operator : CG/JU
 Sample : L4359-10DL 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CB7P1DL

Manual Integrations
 APPROVED

mohammad
 10/15/2020 8:20:20 AM

Quant Time: Oct 15 03:38:45 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 15 00:34:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	66010	20.00	ng/ul	0.00
18) Naphthalene-d8	10.90	136	254086	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.72	164	189155	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.46	188	469742	20.00	ng/ul	0.00
78) Chrysene-d12	21.73	240	547557	20.00	ng/ul	0.00
86) Perylene-d12	25.00	264	554238	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	7.24	99	5105m	0.90	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.41	67	10151	2.95	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	11654	2.83	ng/ul	0.00
13) 4-Methylphenol-d8	8.79	113	8891	2.01	ng/ul	0.00
19) Nitrobenzene-d5	9.25	128	7188	3.78	ng/ul	0.00
22) 2-Nitrophenol-d4	9.98	143	8063	4.17	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.52	165	15348	3.38	ng/ul	0.00
29) 4-Chloroaniline-d4	11.03	131	19915	3.69	ng/ul	0.00
44) Dimethylphthalate-d6	14.12	166	57461	3.98	ng/ul	0.00
47) Acenaphthylene-d8	14.41	160	59567	3.47	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.71	176	50975	3.83	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.82	200	5034	2.14	ng/ul	0.00
71) Anthracene-d10	17.56	188	86446m	4.01	ng/ul	0.00
79) Pyrene-d10	19.84	212	108426	3.82	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.77	264	111458	3.73	ng/ul	0.00

Target Compounds

				Ovalue		
2) 1,4-Dioxane	3.55	88	17395	10.090	ng/uL#	97
45) Dimethylphthalate	14.17	163	25377	1.735	ng/ul#	87

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

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