

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP090919\
 Data File : BP000138.D
 Acq On : 09 Sep 2019 17:08
 Operator : HP/JU
 Sample : K4708-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB5Q6

Quant Time: Sep 10 03:36:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 02:43:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	451812	20.00	ng/ul	0.00
18) Naphthalene-d8	8.18	136	1766474	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	989578	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1832873	20.00	ng/ul	0.00
78) Chrysene-d12	14.13	240	1569036	20.00	ng/ul	-0.01
86) Perylene-d12	15.67	264	1588195	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.60	96	66845	5.25	ng/uL	0.00
5) Phenol-d5	6.51	99	1110839	28.87	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.58	67	575917	28.62	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	900271	28.93	ng/ul	0.00
13) 4-Methylphenol-d8	7.26	113	848335	29.34	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	421798	31.45	ng/ul	0.00
22) 2-Nitrophenol-d4	7.78	143	443644	33.97	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.02	165	784549	28.31	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	1111366	32.43	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	2141149	29.30	ng/ul	0.00
47) Acenaphthylene-d8	9.79	160	2637288	29.21	ng/ul	0.00
52) 4-Nitrophenol-d4	10.03	143	361297	27.67	ng/ul	0.00
58) Fluorene-d10	10.47	176	1776794	28.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.52	200	213557	25.35	ng/ul	0.00
71) Anthracene-d10	11.50	188	2535308	28.56	ng/ul	0.00
79) Pyrene-d10	12.89	212	2639351	28.28	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.58	264	2428079	29.50	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.52	94	87933	2.192	ng/ul 99
45) Dimethylphthalate	9.65	163	619333	8.399	ng/ul 99

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

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