

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
 Data File : BM022591.D
 Acq On : 09 Sep 2019 06:56
 Operator : HP/JU
 Sample : K4706-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BF567

Quant Time: Sep 09 07:47:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 09 07:45:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	372770	20.00	ng/ul	0.00
18) Naphthalene-d8	10.65	136	1671408	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.49	164	1065275	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.22	188	2213400	20.00	ng/ul	0.00
78) Chrysene-d12	21.39	240	2230503	20.00	ng/ul	0.00
86) Perylene-d12	23.68	264	2591221	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.30	96	46564	5.20	ng/uL	0.00
5) Phenol-d5	7.01	99	241985	7.04	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.18	67	464376	23.56	ng/ul	0.00
9) 2-Chlorophenol-d4	7.39	132	553215	20.37	ng/ul	0.00
13) 4-Methylphenol-d8	8.55	113	415364	14.86	ng/ul	0.00
19) Nitrobenzene-d5	9.00	128	304763	25.27	ng/ul	0.00
22) 2-Nitrophenol-d4	9.73	143	293318	26.38	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.26	165	642583	22.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.77	131	63428	1.70	ng/ul	0.00
44) Dimethylphthalate-d6	13.89	166	2230808	25.84	ng/ul	0.00
47) Acenaphthylene-d8	14.18	160	2537833	24.55	ng/ul	0.00
52) 4-Nitrophenol-d4	14.66	143	94109	6.38	ng/ul	0.00
58) Fluorene-d10	15.47	176	1896280	26.24	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.58	200	228691	22.73	ng/ul	0.00
71) Anthracene-d10	17.32	188	3051511	26.91	ng/ul	0.00
79) Pyrene-d10	19.61	212	3444322	27.83	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.54	264	3789316	27.21	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.03	94	73068	2.034 ng/ul	99
45) Dimethylphthalate	13.94	163	225929	2.578 ng/ul	99

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

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