

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\
 Data File : BM023768.D
 Acq On : 16 Nov 2019 00:13
 Operator : JU
 Sample : K5699-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F2J03

Quant Time: Nov 16 03:21:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 16 03:18:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	308001	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1271607	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.18	164	809911	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1799801	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1839675	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	2194324	20.00	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.06	96	52344	7.01	ng/uL	0.00
5) Phenol-d5	6.72	99	207670	7.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	511761	31.72	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	580170	28.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.24	113	383009	17.10	ng/ul	0.00
19) Nitrobenzene-d5	8.67	128	334697	36.77	ng/ul	0.00
22) 2-Nitrophenol-d4	9.39	143	373165	39.01	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	730149	33.47	ng/ul	0.00
29) 4-Chloroaniline-d4	10.45	131	32703	1.37	ng/ul	0.00
44) Dimethylphthalate-d6	13.60	166	2312285	36.61	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	2572822	35.69	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	69709	6.47	ng/ul	0.12
58) Fluorene-d10	15.18	176	1960375	35.15	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	337117	32.38	ng/ul	0.00
71) Anthracene-d10	17.03	188	3207432	37.57	ng/ul	0.00
79) Pyrene-d10	19.33	212	3728714	38.87	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.15	264	4466257	38.68	ng/ul	-0.01

Target Compounds					Ovalue	
6) Phenol	6.74	94	36211	1.252	ng/ul	96
45) Dimethylphthalate	13.65	163	394936	6.381	ng/ul	99

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

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