

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\
 Data File : BM023767.D
 Acq On : 15 Nov 2019 23:37
 Operator : JU
 Sample : K5699-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F2J04

Quant Time: Nov 16 03:21:52 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	310057	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1278142	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.18	164	799646	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1704674	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1683146	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	2016835	20.00	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.06	96	60641	8.06	ng/uL	0.00
5) Phenol-d5	6.72	99	139327	4.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	395565	24.36	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	429785	20.99	ng/ul	0.00
13) 4-Methylphenol-d8	8.24	113	264978	11.75	ng/ul	0.00
19) Nitrobenzene-d5	8.67	128	262222	28.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.39	143	296845	30.87	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	546098	24.91	ng/ul	0.00
29) 4-Chloroaniline-d4	10.46	131	8011	0.33	ng/ul	0.01
44) Dimethylphthalate-d6	13.60	166	1906140	30.57	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	2100764	29.52	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	40286	3.78	ng/ul	0.00
58) Fluorene-d10	15.18	176	1620284	29.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	262362	26.61	ng/ul	0.00
71) Anthracene-d10	17.03	188	2589092	32.02	ng/ul	0.00
79) Pyrene-d10	19.33	212	2969007	33.83	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.15	264	3526082	33.22	ng/ul	-0.01

Target Compounds					Ovalue	
6) Phenol	6.74	94	36715	1.261	ng/ul	98
45) Dimethylphthalate	13.65	163	365262	5.977	ng/ul	99

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

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