

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111819\
 Data File : BM023779.D
 Acq On : 18 Nov 2019 13:18
 Operator : JU
 Sample : K5699-15
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F2J13

Quant Time: Nov 18 14:01:13 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	302513	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1239040	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.18	164	768518	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1633709	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1640323	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	1980048	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	47322	6.45	ng/uL	0.00
5) Phenol-d5	6.71	99	160568	5.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	448130	28.28	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	469918	23.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.24	113	305083	13.87	ng/ul	0.00
19) Nitrobenzene-d5	8.68	128	282336	31.83	ng/ul	0.00
22) 2-Nitrophenol-d4	9.39	143	308942	33.14	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	576850	27.14	ng/ul	0.00
29) 4-Chloroaniline-d4	10.45	131	27978	1.20	ng/ul	0.00
44) Dimethylphthalate-d6	13.60	166	1974712	32.95	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	2166991	31.68	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	39549	3.87	ng/ul	0.00
58) Fluorene-d10	15.18	176	1653031	31.23	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	272428	28.83	ng/ul	0.00
71) Anthracene-d10	17.03	188	2563407	33.08	ng/ul	0.00
79) Pyrene-d10	19.33	212	3049793	35.66	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.15	264	3672572	35.25	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
45) Dimethylphthalate	13.65	163	164216	2.796	ng/ul	98

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

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