

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111819\
 Data File : BM023778.D
 Acq On : 18 Nov 2019 12:41
 Operator : JU
 Sample : K5699-14
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F2J12

Quant Time: Nov 18 13:58:02 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	297430	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1218296	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.18	164	754657	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1610940	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1613064	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	1956954	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	46906	6.50	ng/uL	0.00
5) Phenol-d5	6.71	99	190744	7.04	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	490586	31.49	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	522974	26.63	ng/ul	0.00
13) 4-Methylphenol-d8	8.24	113	342491	15.84	ng/ul	0.00
19) Nitrobenzene-d5	8.67	128	311261	35.69	ng/ul	0.00
22) 2-Nitrophenol-d4	9.39	143	343116	37.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	607132	29.05	ng/ul	0.00
29) 4-Chloroaniline-d4	10.45	131	56364	2.46	ng/ul	0.00
44) Dimethylphthalate-d6	13.60	166	2032049	34.53	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	2315054	34.47	ng/ul	0.00
52) 4-Nitrophenol-d4	14.41	143	52118	5.19	ng/ul	0.00
58) Fluorene-d10	15.18	176	1800100	34.64	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	294094	31.56	ng/ul	0.00
71) Anthracene-d10	17.03	188	2979745	38.99	ng/ul	0.00
79) Pyrene-d10	19.33	212	3356031	39.90	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.15	264	4069679	39.52	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
45) Dimethylphthalate	13.65	163	137727	2.388	ng/ul	99

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

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