

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\
 Data File : BM023765.D
 Acq On : 15 Nov 2019 22:25
 Operator : JU
 Sample : K5699-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F2J01

Quant Time: Nov 16 04:15:28 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	309112	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1241865	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.18	164	775782	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1672812	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1654415	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	1947259	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	45360	6.05	ng/uL	0.00
5) Phenol-d5	6.72	99	160520	5.70	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	421972	26.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.06	132	466990	22.88	ng/ul	0.00
13) 4-Methylphenol-d8	8.24	113	295614	13.15	ng/ul	0.00
19) Nitrobenzene-d5	8.67	128	275440	30.99	ng/ul	0.00
22) 2-Nitrophenol-d4	9.39	143	305405	32.69	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.93	165	587380	27.57	ng/ul	0.00
29) 4-Chloroaniline-d4	10.45	131	16323	0.70	ng/ul	0.00
44) Dimethylphthalate-d6	13.60	166	1948441	32.21	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	2182160	31.61	ng/ul	0.00
52) 4-Nitrophenol-d4	14.43	143	47760	4.62	ng/ul	0.02
58) Fluorene-d10	15.18	176	1675277	31.36	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	273136	28.23	ng/ul	0.00
71) Anthracene-d10	17.03	188	2576405	32.47	ng/ul	0.00
79) Pyrene-d10	19.33	212	2997031	34.74	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	3523613	34.39	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.74	94	82463	2.841	ng/ul 98
28) Naphthalene	10.35	128	272578	4.176	ng/ul 99
45) Dimethylphthalate	13.65	163	318782	5.377	ng/ul 99

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

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