

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
 Data File : BM023945.D
 Acq On : 28 Nov 2019 10:42
 Operator : JU
 Sample : K5915-06
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLNW1

Quant Time: Nov 28 21:36:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 27 17:16:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	329916	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	1323813	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.15	164	790461	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	1658494	20.00	ng/ul	0.00
78) Chrysene-d12	21.11	240	1702476	20.00	ng/ul	0.00
86) Perylene-d12	23.25	264	2090591	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.03	96	37399	5.04	ng/uL	0.00
5) Phenol-d5	6.67	99	664916	22.99	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	435924	25.92	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	556897	24.92	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	395292	17.33	ng/ul	0.00
19) Nitrobenzene-d5	8.64	128	270934	26.37	ng/ul	0.00
22) 2-Nitrophenol-d4	9.36	143	304264	26.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	548314	24.79	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	581473	23.96	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	1757612	28.62	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	2092997	29.08	ng/ul	0.00
52) 4-Nitrophenol-d4	14.38	143	232604	22.53	ng/ul	0.00
58) Fluorene-d10	15.15	176	1556592	30.12	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	211415	20.31	ng/ul	0.00
71) Anthracene-d10	17.00	188	2365682	29.87	ng/ul	0.00
79) Pyrene-d10	19.30	212	2827271	32.14	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.12	264	3443840	31.16	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.70	94	39722	1.335	ng/ul 93
45) Dimethylphthalate	13.62	163	634966	10.561	ng/ul 100

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

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